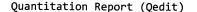
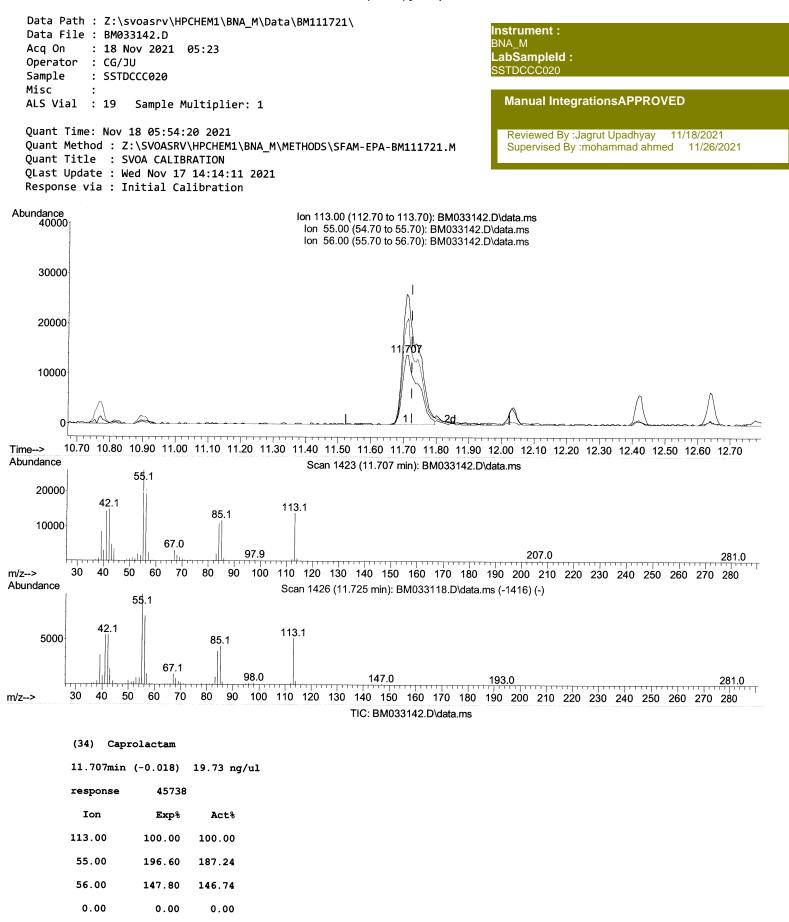
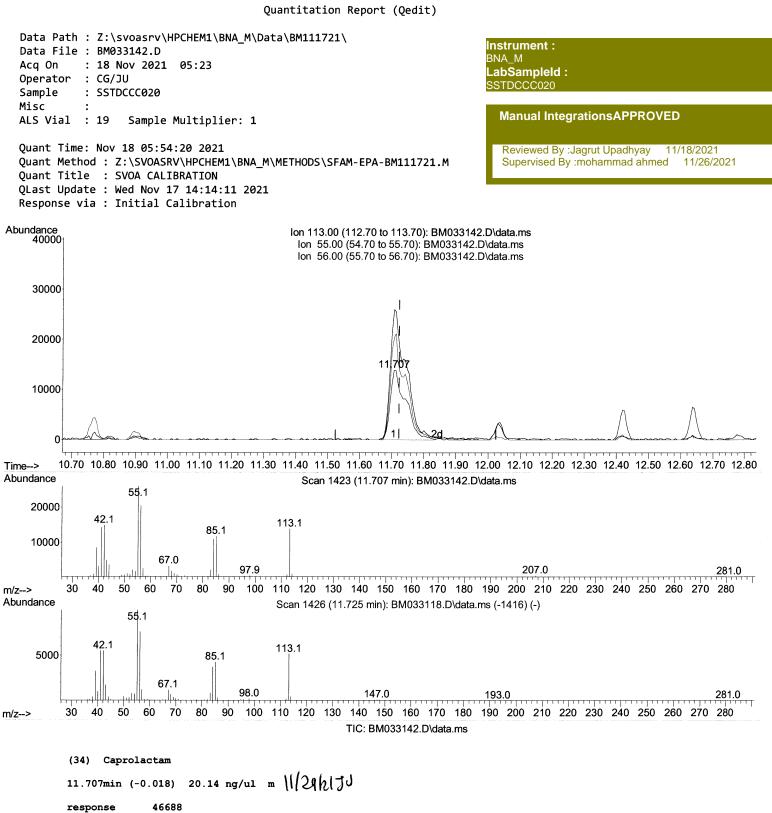


SFAM-EPA-BM111721.M Thu Nov 18 06:12:42 2021







response	40000		
Ion	Exp%	Act%	
113.00	100.00	100.00	
55.00	196.60	187.24	
56.00	147.80	146.74	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\H Data File : BM033142.D Acq On : 18 Nov 2021 Operator : CG/JU Sample : SSTDCCC020 Misc : ALS Vial : 19 Sample M	_	Instrument : BNA_M LabSampleId : SSTDCCC020 Manual IntegrationsAPPROVED			
Quant Time: Nov 18 05:54 Quant Method : Z:\SVOASR Quant Title : SVOA CALIE QLast Update : Wed Nov 17 Response via : Initial Ca	/\HPCHEM1\BNA_M 3RATION 7 14:14:11 2021	Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021			
Compound		-	•	Conc Units Dev	
Internal Standards					
1) 1,4-Dichlorobenzene	e-d4 7.972	152	124596	20.000 ng/ul	0.00
20) Naphthalene-d8	10.766		504985	20.000 ng/ul	-0.01
38) Acenaphthene-d10	14.589		348889	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.324		781815	20.000 ng/ul	0.00
79) Chrysene-d12	21.477		759707	20.000 ng/ul	0.00
88) Perylene-d12	23.824		747245	20.000 ng/ul	-0.01
· ·				8,	
System Monitoring Compou	inds				
3) 1,4-Dioxane-d8	3.419	96	22715	7.028 ng/uL	0.00
4) Pyridine-d5	3.843	84	151299	16.995 ng/ul	0.00
7) Phenol-d5	7.125	99	184067	17.465 ng/ul	0.00
9) Bis-(2-Chloroethyl)	eth 7.301	67	118982	17.799 ng/ul	0.00
11) 2-Chlorophenol-d4	7.501	132	143048	17.920 ng/ul	0.00
15) 4-Methylphenol-d8	8.666	113	146464	17.907 ng/ul	0.00
21) Nitrobenzene-d5	9.131	128	69188	19.083 ng/ul	0.00
24) 2-Nitrophenol-d4	9.848	143	71394	19.662 ng/ul	-0.01
28) 2,4-Dichlorophenol-	d3 10.384	165	151754	18.250 ng/ul	0.00
31) 4-Chloroaniline-d4	10.901		199177	18.012 ng/ul	0.00
46) Dimethylphthalate-d			500430	19.555 ng/ul	-0.01
<pre>49) Acenaphthylene-d8</pre>	14.283		588112	17.839 ng/ul	0.00
54) 4-Nitrophenol-d4	14.772		81253	19.492 ng/ul	0.00
60) Fluorene-d10	15.577		435971	18.991 ng/ul	0.00
65) 4,6-Dinitro-2-methy			70663	19.232 ng/ul	0.00
73) Anthracene-d10	17.424		695088	18.444 ng/ul	0.00
81) Pyrene-d10	19.701		798622	17.789 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.671	264	730845	18.224 ng/ul	-0.01
Target Compounds				Ova	alue
2) 1,4-Dioxane	3.455	88	22983	7.004 ng/uL#	
5) Pyridine	3.860	79	157381	17.324 ng/ul	99
6) Benzaldehyde	7.113	77	135494	22.558 ng/ul	98
8) Phenol	7.149	94	184214	17.552 ng/ul	96
Bis(2-Chloroethyl)e	ther 7.390	93	151369	18.156 ng/ul	98
12) 2-Chlorophenol	7.531	128	147837	17.943 ng/ul	98
13) 2-Methylphenol	8.401	108	141189	17.561 ng/ul	95
14) 2,2'-oxybis(1-Chlor		45	239576	18.606 ng/ul	100
16) Acetophenone	8.795	105	235075	18.209 ng/ul	96
17) N-Nitroso-di-n-prop		70	128134	18.421 ng/ul	98
18) 4-Methylphenol	8.731	108	151378	17.961 ng/ul	98
19) Hexachloroethane	9.048	117	66676	17.764 ng/ul	94
22) Nitrobenzene	9.172	77	187374	18.519 ng/ul	99
23) Isophorone 25) 2-Nitrophenol	9.701 9.884	82 139	341281 76300	18.454 ng/ul 19.824 ng/ul	99 90
26) 2,4-Dimethylphenol	9.931	107	180309	17.830 ng/ul	96
27) Bis(2-Chloroethoxy)		93	201942	18.349 ng/ul	98
29) 2,4-Dichlorophenol	10.407	162	148494	18.416 ng/ul	98
30) Naphthalene	10.819	128	483658	18.011 ng/ul	99
32) 4-Chloroaniline	10.925	127	485058 197491	17.679 ng/ul	96
33) Hexachlorobutadiene	11.101	225	109146	17.378 ng/ul	98
34) Caprolactam	11.707	113		20.141 ng/ul >	
35) 4-Chloro-3-methylph			172662	19.665 ng/ul	99

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	Quantitation Report (QT Reviewed)										
Data Path : Z:\svoasrv\HPCHEM1\	BNA M\Dat	a\BM1	11721\								
Data File : BM033142.D					Instrument :						
Acq On : 18 Nov 2021 05:23					BNA_M						
Operator : CG/JU					LabSampleId : SSTDCCC020						
Sample : SSTDCCC020					5515666020						
Misc :					Manual IntegrationsAPPROVED						
ALS Vial : 19 Sample Multipl	ier: 1										
Quant Time: New 18 AF. F4.20 202											
Quant Time: Nov 18 05:54:20 202 Quant Method : Z:\SVOASRV\HPCHE	Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021										
Quant Title : SVOA CALIBRATION	T (DINA_III)			A-DHIII/21.H	Supervised by Inchammad anned 11/20/2021						
QLast Update : Wed Nov 17 14:14	:11 2021										
Response via : Initial Calibration											
Compound	R.T.	QIon	Response	Conc Units Dev(Min)						
26 2 Mothylpephthelene	12 410	140									
36) 2-Methylnaphthalene 37) 1-Methylnaphthalene	12.419 12.636	142 142	338491 345097	18.193 ng/ul 18.169 ng/ul	97 98						
39) 1,2,4,5-Tetrachloroben	12.783		196749	16.950 ng/ul	97						
40) Hexachlorocyclopentadiene		237	132108	16.175 ng/ul	96						
41) 2,4,6-Trichlorophenol	13.019	196	124968	18.281 ng/ul	97						
42) 2,4,5-Trichlorophenol	13.089	196	140413	19.155 ng/ul	100						
43) 1,1'-Biphenyl	13.425	154	472755	17.391 ng/ul	99						
44) 2-Chloronaphthalene	13.466	162	357693	17.063 ng/ul	97						
45) 2-Nitroaniline	13.672	65	115895	20.975 ng/ul	100						
47) Dimethylphthalate	14.042	163	485847	19.479 ng/ul	99						
48) 2,6-Dinitrotoluene	14.160	165	90507	21.209 ng/ul	97						
50) Acenaphthylene	14.313	152	589453	17.619 ng/ul	99						
51) 3-Nitroaniline 52) Acenaphthene	14.489 14.654	138 153	95499 390460	22.210 ng/ul 17.891 ng/ul	90 100						
53) 2,4-Dinitrophenol	14.689	184	42639	19.576 ng/ul#	90						
55) 4-Nitrophenol	14.783	109	80344	18.871 ng/ul	97						
56) Dibenzofuran	14.983	168	588119	18.386 ng/ul	98						
57) 2,4-Dinitrotoluene	14.942	165	128514	21.897 ng/ul#	97						
58) 2,3,4,6-Tetrachlorophenol	15.207	232	120443	20.130 ng/ul	98						
59) Diethylphthalate	15.401	149	494899	19.810 ng/ul	99						
61) Fluorene	15.630	166	481604	18.943 ng/ul	98						
62) 4-Chlorophenyl-phenyle	15.624	204	249420	18.796 ng/ul	99						
63) 4-Nitroaniline	15.648	138	96574	22.842 ng/ul	95						
66) 4,6-Dinitro-2-methylph 67) N-Nitrosodiphenylamine	15.701 15.836	198 169	71240 423627	19.345 ng/ul	99 98						
68) 4-Bromophenyl-phenylether		248	423627	18.423 ng/ul 18.020 ng/ul	98 97						
69) Hexachlorobenzene	16.624	284	177109	17.782 ng/ul	97						
70) Atrazine	16.783	200	158932	17.887 ng/ul	98						
71) Pentachlorophenol	16.966	266	105117	18.258 ng/ul	99						
72) Phenanthrene	17.365	178	788410	18.205 ng/ul	99						
74) Anthracene	17.454	178	795627	18.320 ng/ul	98						
75) 1,2,3,4-Tetrachloroben	13.389	216	203426	15.841 ng/uL	97						
76) Pentachlorobenzene	14.901	250	218887	17.387 ng/uL	99						
77) Carbazole	17.724	167	704506	18.260 ng/ul	99						
78) Di-n-butylphthalate 80) Fluoranthene	18.283 19.371	149 202	829873 937040	19.671 ng/ul 17.816 ng/ul	100 99						
82) Pyrene	19.730	202	962109	18.016 ng/ul	99						
83) Butylbenzylphthalate	20.618	149	352035	19.295 ng/ul	98						
84) 3,3'-Dichlorobenzidine	21.395	252	313134	18.296 ng/ul	100						
85) Benzo(a)anthracene	21.459	228	897684	18.269 ng/ul	99						
<pre>86) Bis(2-ethylhexyl)phtha</pre>	21.383	149	503204	19.458 ng/ul	98						
87) Chrysene	21.512	228	869806	18.127 ng/ul	99						
89) Di-n-octyl phthalate		149	839385	18.436 ng/ul	100						
90) Benzo(b)fluoranthene			928675	18.453 ng/ul	99						
91) Benzo(k)fluoranthene		252	831769	18.042 ng/ul	99						
93) Benzo(a)pyrene 94) Indeno(1,2,3-cd)pyrene	23.718 26.224	252 276	873908 955676	18.374 ng/ul 18.044 ng/ul	98 96						
95) Dibenzo(a,h)anthracene	26.224		955676 818523	18.064 ng/ul	96 98						
96) Benzo(g,h,i)perylene	26.965		825609	17.960 ng/ul	99						

(#) = qualifier out of range (m) = manual integration (+) = signals summed