Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112221\

Data File : BN017542.D

Acq On : 22 Nov 2021 12:14

Operator : CG/JU Sample : SSTD04048

Misc :

ALS Vial : 5 Sample Multiplier: 1

Quant Time: Nov 22 14:52:41 2021

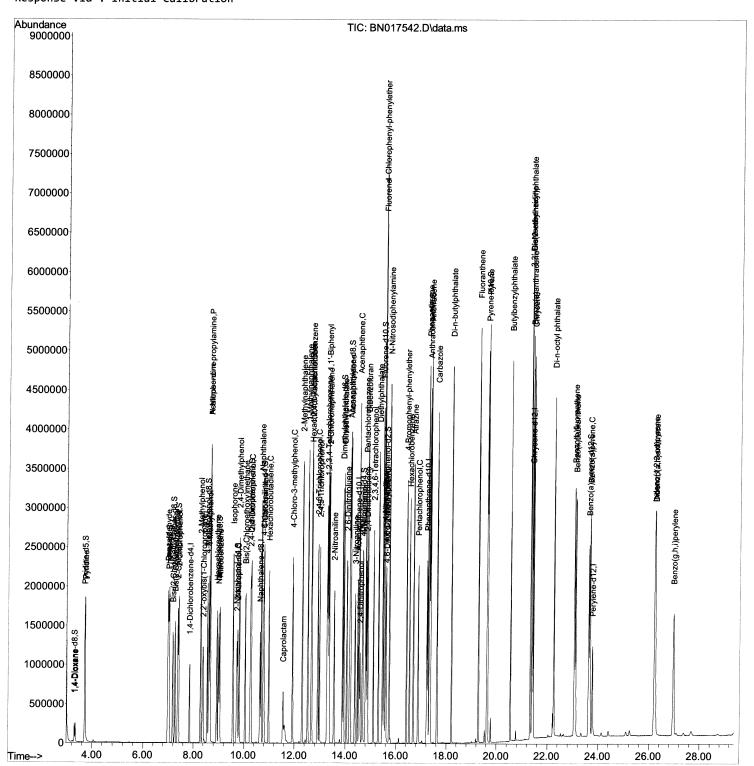
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION
OLast Undate : Mon Nov 22 14:49:39 20

QLast Update : Mon Nov 22 14:49:39 2021 Response via : Initial Calibration

Instrument : BNA_N ClientSampleId : SSTD040248

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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Data File : BN017542.D

Acq On : 22 Nov 2021 12:14

Operator : CG/JU Sample : SSTD04048

Misc

ALS Vial : 5 Sample Multiplier: 1

Quant Time: Nov 22 14:52:41 2021

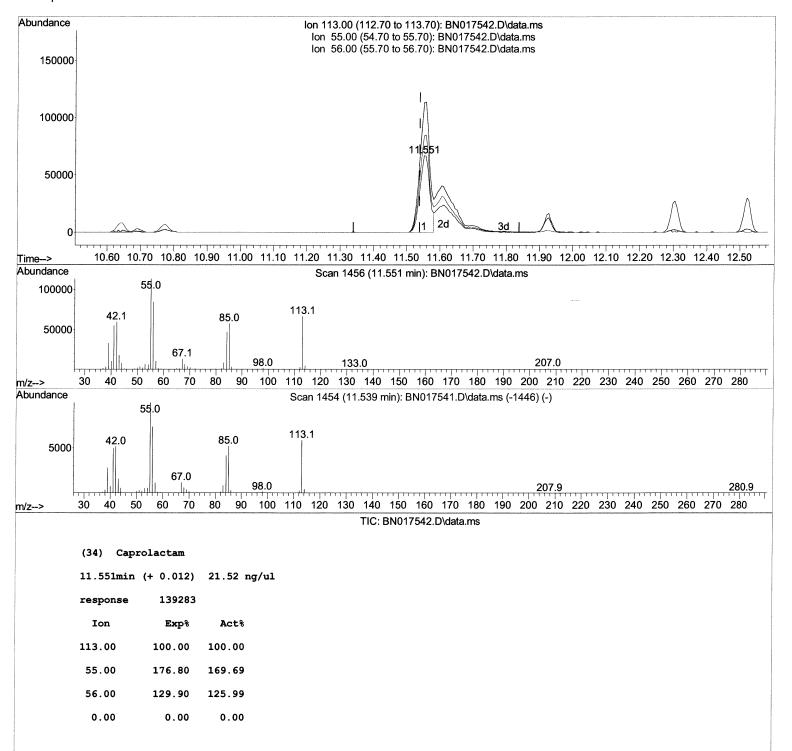
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

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Misc

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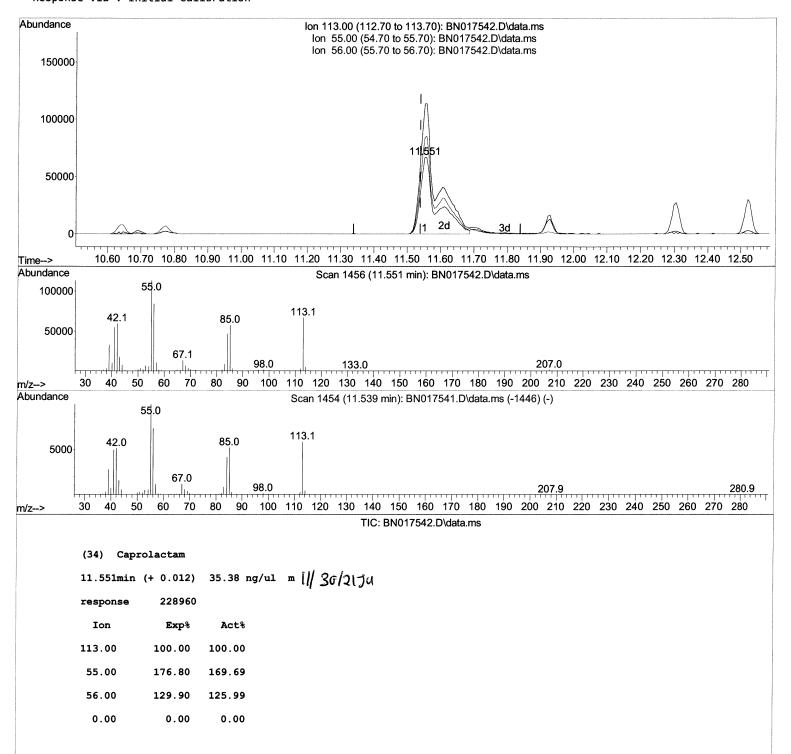
Quant Time: Nov 22 14:52:41 2021

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BNA_N
ClientSampleId:
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Manual IntegrationsAPPROVED



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 $\label{thm:local_NMethods} Quant \ \mbox{Methods} : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M$

Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 14:49:39 2021 Response via : Initial Calibration Instrument: BNA_N ClientSampleId: SSTD040248

Manual IntegrationsAPPROVED

Compound			Response			
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	260548	20.000	ng/ul	0.00
20) Naphthalene-d8	10.640	136	1152989		ng/ul	0.00
38) Acenaphthene-d10	14.481		702697		ng/ul	0.00
64) Phenanthrene-d10	17.222		1362758		ng/ul	0.00
79) Chrysene-d12	21.410		915799		ng/ul	0.00
88) Perylene-d12	23.774		778582		ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	115073		ng/uL	0.00
4) Pyridine-d5	3.705	84	795331		ng/ul	0.00
7) Phenol-d5	7.011	99	986621		ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.175	67	607576		ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.375	132	758796		ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.552	113	786640		ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	379645	42.517	_	0.00
24) 2-Nitrophenol-d4	9.722	143	390344	39.487		0.00
28) 2,4-Dichlorophenol-d3	10.263	165	720055	39.759	-	0.00
31) 4-Chloroaniline-d4	10.775	131	1032926	38.294		0.00
46) Dimethylphthalate-d6	13.892	166	2116694	40.237		0.00
<pre>49) Acenaphthylene-d8</pre>	14.175	160	2688230	40.537		0.00
54) 4-Nitrophenol-d4	14.675	143	376571	36.279		0.01
60) Fluorene-d10	15.475	176	1688539	37.580		0.00
65) 4,6-Dinitro-2-methylph	15.592	200	296816	33.636		0.00
73) Anthracene-d10	17.322	188	2545811	38.693	-	0.00
81) Pyrene-d10	19.616	212	2592086	48.261		0.00
92) Benzo(a)pyrene-d12	23.627	264	1732807	41.876	ng/ul	0.00
Target Compounds					Qva:	lue
2) 1,4-Dioxane	3.340	88	116578	16.389	•	95
5) Pyridine	3.728	79	796394	40.847	-	91
6) Benzaldehyde	6.981	77	627179	43.477	•	92
8) Phenol	7.034	94	997279	40.718		93
10) Bis(2-Chloroethyl)ether	7.269	93	816847	42.444		98
12) 2-Chlorophenol	7.410	128	775352	39.379		97
13) 2-Methylphenol	8.287	108	763699	40.984		98
14) 2,2'-oxybis(1-Chloropr	8.381	45	1095045	38.785		96
16) Acetophenone	8.663	105	1167641	39.863	-	93
17) N-Nitroso-di-n-propyla	8.663	70	621412	42.261		94
18) 4-Methylphenol	8.616	108	816092	39.642	-	99
19) Hexachloroethane	8.928	117	322838	42.575		94
22) Nitrobenzene	9.040	77	922250	45.931		94
23) Isophorone	9.569	82	1777532	43.602		96
25) 2-Nitrophenol	9.752	139	420766		ng/ul#	89
26) 2,4-Dimethylphenol	9.816	107	909803	42.179		94
27) Bis(2-Chloroethoxy)met	10.057	93	1096799	42.063		98
29) 2,4-Dichlorophenol	10.287	162	704634	39.175		98
30) Naphthalene	10.693	128	2470141	39.377	_	100
32) 4-Chloroaniline	10.799	127	1044903	39.158		99
33) Hexachlorobutadiene	10.993	225	458824	41.891		95
34) Caprolactam	11.551	113	228960m >			1113012174
35) 4-Chloro-3-methylphenol	11.928	107	823643	41.157		95
					-	

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Misc

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Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 14:49:39 2021 Response via : Initial Calibration

Instrument : BNA_N ClientSampleId : SSTD040248

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.304	142	1652959	38.167 ng/ul	100
37) 1-Methylnaphthalene	12.522	142	1668961	37.410 ng/ul	97
39) 1,2,4,5-Tetrachloroben	12.675	216	821190	40.094 ng/ul#	95
40) Hexachlorocyclopentadiene	12.663		525460	40.367 ng/ul	96
41) 2,4,6-Trichlorophenol	12.910	196	550254	41.435 ng/ul	98
42) 2,4,5-Trichlorophenol	12.981	196	590900	39.957 ng/ul	98
43) 1,1'-Biphenyl	13.316	154	2181482	40.096 ng/ul	100
44) 2-Chloronaphthalene	13.357	162	1666494	40.219 ng/ul	95
45) 2-Nitroaniline	13.551	65	533516	47.300 ng/ul	92
47) Dimethylphthalate	13.940	163	2060059	38.992 ng/ul	99
48) 2,6-Dinitrotoluene	14.051	165	434448	41.877 ng/ul	91
50) Acenaphthylene	14.204	152	2668150	39.284 ng/ul	100
51) 3-Nitroaniline	14.381	138	451972	37.640 ng/ul	95
52) Acenaphthene	14.545	153	1712301	38.986 ng/ul	99
53) 2,4-Dinitrophenol	14.581	184	205665	30.995 ng/ul	93
55) 4-Nitrophenol	14.687	109	324029	44.237 ng/ul#	85
56) Dibenzofuran	14.881	168	2427984	38.719 ng/ul	97
57) 2,4-Dinitrotoluene	14.839	165	599631	40.078 ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	15.104	232	436738	35.775 ng/ul	96
59) Diethylphthalate	15.310	149	2120436	40.249 ng/ul	98
61) Fluorene	15.528	166	1759174	35.590 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.522	204	871053	35.838 ng/ul	90
63) 4-Nitroaniline	15.545	138	396281	33.399 ng/ul#	87
66) 4,6-Dinitro-2-methylph	15.604	198	295329	33.544 ng/ul#	88
67) N-Nitrosodiphenylamine	15.734	169	1651704	39.820 ng/ul	98
68) 4-Bromophenyl-phenylether	16.416	248	541776	37.307 ng/ul#	89
69) Hexachlorobenzene	16.539	284	573488	33.936 ng/ul	92
70) Atrazine	16.686	200	624018	41.000 ng/ul	97
71) Pentachlorophenol	16.875	266	347767	34.279 ng/ul	94
72) Phenanthrene	17.269	178	2905107	38.691 ng/ul	98
74) Anthracene	17.357	178	2883807	37.927 ng/ul	97
75) 1,2,3,4-Tetrachloroben	13.281	216	856771	43.008 ng/uL	97
76) Pentachlorobenzene	14.804	250	771991	38.267 ng/uL	99
77) Carbazole	17.622	167	2599555	38.723 ng/ul	99
78) Di-n-butylphthalate	18.198	149	3348600	42.199 ng/ul	99
80) Fluoranthene	19.280	202	3149035	49.098 ng/ul	100
82) Pyrene	19.645	202	3066829	46.714 ng/ul	99
83) Butylbenzylphthalate	20.551	149	1266832	51.825 ng/ul	92
84) 3,3'-Dichlorobenzidine	21.327	252	748503	36.716 ng/ul	95
85) Benzo(a)anthracene	21.392	228	2503702	41.530 ng/ul	97
86) Bis(2-ethylhexyl)phtha	21.339	149	1731929	46.944 ng/ul#	99
87) Chrysene 89) Di-n-octyl phthalate	21.445	228	2422467	40.769 ng/ul	99
- ·	22.251	149	2701896	47.480 ng/ul	100
90) Benzo(b)fluoranthene91) Benzo(k)fluoranthene	23.063	252	2211090	42.066 ng/ul	95
	23.110 23.674	252	2134138	41.741 ng/ul	98
93) Benzo(a)pyrene 94) Indeno(1,2,3-cd)pyrene	26.221	252 276	2106534 2238450	41.650 ng/ul 43.385 ng/ul	99 99
95) Dibenzo(a,h)anthracene	26.245	278	1913901	43.884 ng/ul	99 97
96) Benzo(g,h,i)perylene	26.243	276	1925336	44.567 ng/ul	100
				IIB/UI	

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