

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036582.D
 Acq On : 19 Sep 2022 16:21
 Operator : CG/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD020177

Quant Time: Sep 19 22:32:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 09/20/2022
 Supervised By :mohammad ahmed 09/21/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	345556	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	1526983	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	972717	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	1934550	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	1894580	20.000	ng/u1	0.00
88) Perylene-d12	23.997	264	1786644	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	69252	7.513	ng/uL	0.00
4) Pyridine-d5	3.816	84	508797	18.926	ng/u1	0.00
7) Phenol-d5	7.175	99	579293	18.759	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	363818	19.400	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	443815	19.962	ng/u1	0.00
15) 4-Methylphenol-d8	8.734	113	451082	19.769	ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	216718	20.969	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	197937	21.381	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	430912	20.085	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	688017	19.288	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	1276924	19.464	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	1519799	19.530	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	238264	18.725	ng/u1	0.00
60) Fluorene-d10	15.639	176	1141320	19.413	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	144216	18.842	ng/u1	0.00
73) Anthracene-d10	17.486	188	1720341	19.449	ng/u1	0.00
81) Pyrene-d10	19.768	212	1880026	20.171	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.838	264	1728931	19.807	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.428	88	74000	7.452	ng/uL	99
5) Pyridine	3.834	79	496452	18.381	ng/u1	99
6) Benzaldehyde	7.157	77	323837m	21.656	ng/u1	
8) Phenol	7.204	94	626274	18.896	ng/u1	99
10) Bis(2-Chloroethyl)ether	7.445	93	503369	19.520	ng/u1	100
12) 2-Chlorophenol	7.587	128	484118	20.130	ng/u1	97
13) 2-Methylphenol	8.463	108	471494	19.354	ng/u1	99
14) 2,2'-oxybis(1-Chloropr...	8.563	45	602134	19.701	ng/u1	99
16) Acetophenone	8.851	105	791595	20.001	ng/u1	99
17) N-Nitroso-di-n-propyla...	8.839	70	395752	20.723	ng/u1	98
18) 4-Methylphenol	8.792	108	521057	19.780	ng/u1	100
19) Hexachloroethane	9.116	117	194454	19.962	ng/u1	94
22) Nitrobenzene	9.233	77	566604	20.648	ng/u1	98
23) Isophorone	9.763	82	1114963	19.927	ng/u1	100
25) 2-Nitrophenol	9.945	139	237637	20.783	ng/u1	95
26) 2,4-Dimethylphenol	10.004	107	569796	20.204	ng/u1	99
27) Bis(2-Chloroethoxy)met...	10.245	93	672365	20.161	ng/u1	100
29) 2,4-Dichlorophenol	10.480	162	452962	20.083	ng/u1	100
30) Naphthalene	10.892	128	1646208	19.790	ng/u1	100
32) 4-Chloroaniline	10.998	127	707390	19.208	ng/u1	98
33) Hexachlorobutadiene	11.180	225	269301	19.522	ng/u1	99
34) Caprolactam	11.751	113	139570m	19.076	ng/u1	
35) 4-Chloro-3-methylphenol	12.110	107	509617	20.468	ng/u1	99
36) 2-Methylnaphthalene	12.492	142	1099072	19.831	ng/u1	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036582.D
 Acq On : 19 Sep 2022 16:21
 Operator : CG/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD020177

Quant Time: Sep 19 22:32:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 09/20/2022
 Supervised By :mohammad ahmed 09/21/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.710	142	1100784	19.757	ng/u1	100
39) 1,2,4,5-Tetrachloroben...	12.857	216	535368	19.445	ng/u1	100
40) Hexachlorocyclopentadiene	12.839	237	280045	19.518	ng/u1	99
41) 2,4,6-Trichlorophenol	13.086	196	332804	19.742	ng/u1	99
42) 2,4,5-Trichlorophenol	13.157	196	368683	20.399	ng/u1	98
43) 1,1'-Biphenyl	13.492	154	1424410	19.865	ng/u1	99
44) 2-Chloronaphthalene	13.533	162	1102135	19.588	ng/u1	99
45) 2-Nitroaniline	13.733	65	303608	21.922	ng/u1	99
47) Dimethylphthalate	14.110	163	1362899	19.584	ng/u1	99
48) 2,6-Dinitrotoluene	14.221	165	239146	21.137	ng/u1	99
50) Acenaphthylene	14.374	152	1750369	19.674	ng/u1	100
51) 3-Nitroaniline	14.551	138	269807	19.883	ng/u1	97
52) Acenaphthene	14.715	153	1229817	19.756	ng/u1	99
53) 2,4-Dinitrophenol	14.751	184	98356	19.592	ng/u1	97
55) 4-Nitrophenol	14.839	109	206555	18.989	ng/u1	98
56) Dibenzofuran	15.045	168	1639002	19.605	ng/u1	97
57) 2,4-Dinitrotoluene	15.004	165	352358	21.422	ng/u1	95
58) 2,3,4,6-Tetrachlorophenol	15.268	232	294849	19.489	ng/u1#	99
59) Diethylphthalate	15.468	149	1343135	19.280	ng/u1	100
61) Fluorene	15.692	166	1393635	19.573	ng/u1	99
62) 4-Chlorophenyl-phenyle...	15.686	204	661668	19.431	ng/u1	97
63) 4-Nitroaniline	15.704	138	256255	19.437	ng/u1	100
66) 4,6-Dinitro-2-methylph...	15.762	198	164835	19.085	ng/u1	98
67) N-Nitrosodiphenylamine	15.898	169	1136750	19.951	ng/u1	99
68) 4-Bromophenyl-phenylether	16.580	248	339534	17.807	ng/u1	95
69) Hexachlorobenzene	16.692	284	426324	18.903	ng/u1	98
70) Atrazine	16.845	200	370670	18.436	ng/u1	98
71) Pentachlorophenol	17.033	266	235159	17.554	ng/u1	98
72) Phenanthrene	17.427	178	2113815	19.837	ng/u1	100
74) Anthracene	17.521	178	2117575	19.551	ng/u1	99
75) 1,2,3,4-Tetrachloroben...	13.457	216	527113	19.922	ng/uL	99
76) Pentachlorobenzene	14.968	250	562349	19.650	ng/uL	99
77) Carbazole	17.786	167	1933326	19.280	ng/u1	99
78) Di-n-butylphthalate	18.356	149	2207222	19.420	ng/u1	100
80) Fluoranthene	19.433	202	2331415	20.413	ng/u1	99
82) Pyrene	19.798	202	2480711	20.519	ng/u1	99
83) Butylbenzylphthalate	20.692	149	895019	19.537	ng/u1	93
84) 3,3'-Dichlorobenzidine	21.468	252	748682	20.224	ng/u1	98
85) Benzo(a)anthracene	21.533	228	2366554	19.936	ng/u1	100
86) Bis(2-ethylhexyl)phtha...	21.468	149	1413481	21.122	ng/u1	99
87) Chrysene	21.586	228	2227052	19.765	ng/u1	99
89) Di-n-octyl phthalate	22.415	149	2241562	20.811	ng/u1	100
90) Benzo(b)fluoranthene	23.250	252	2272568	20.058	ng/u1	99
91) Benzo(k)fluoranthene	23.297	252	2268113	20.035	ng/u1	98
93) Benzo(a)pyrene	23.885	252	1972783	19.807	ng/u1	99
94) Indeno(1,2,3-cd)pyrene	26.544	276	2042993	17.565	ng/u1	98
95) Dibenzo(a,h)anthracene	26.562	278	1871467	17.742	ng/u1	99
96) Benzo(g,h,i)perylene	27.326	276	1708506	16.944	ng/u1	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036582.D
 Acq On : 19 Sep 2022 16:21
 Operator : CG/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD020177

Quant Time: Sep 19 22:32:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 09/20/2022
 Supervised By :mohammad ahmed 09/21/2022

