

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092922\
 Data File : BM036796.D
 Acq On : 29 Sep 2022 11:47
 Operator : CG/JU
 Sample : SSTD00578
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD005026

Quant Time: Sep 30 01:11:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 30 01:10:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	325306	20.000	ng/u1	0.00
20) Naphthalene-d8	10.792	136	1479666	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.610	164	974084	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	2001055	20.000	ng/u1	0.00
79) Chrysene-d12	21.515	240	1988440	20.000	ng/u1	0.00
88) Perylene-d12	23.933	264	1990681	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	16624	1.916	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	0.000	99	0d	0.000	ng/u1	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/u1	
11) 2-Chlorophenol-d4	7.510	132	92056	4.398	ng/u1	0.00
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/u1	
21) Nitrobenzene-d5	9.145	128	44295	4.423	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	38631	4.306	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	83670	4.025	ng/u1	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/u1	
46) Dimethylphthalate-d6	14.021	166	292054	4.445	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	306425	3.932	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
60) Fluorene-d10	15.598	176	256651	4.359	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/u1	
73) Anthracene-d10	17.445	188	393097	4.296	ng/u1	0.00
81) Pyrene-d10	19.727	212	428924	4.385	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.774	264	406086	4.175	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.404	88	17406	1.862	ng/uL	95
12) 2-Chlorophenol	7.539	128	102889	4.544	ng/u1	97
17) N-Nitroso-di-n-propyla...	8.792	70	91011	5.062	ng/u1	98
19) Hexachloroethane	9.063	117	42604	4.646	ng/u1	97
22) Nitrobenzene	9.186	77	125716	4.728	ng/u1	98
23) Isophorone	9.722	82	243045	4.483	ng/u1	99
25) 2-Nitrophenol	9.904	139	46503	4.197	ng/u1	93
26) 2,4-Dimethylphenol	9.963	107	120916	4.425	ng/u1	97
27) Bis(2-Chloroethoxy)met...	10.198	93	158299	4.899	ng/u1	98
29) 2,4-Dichlorophenol	10.433	162	90196	4.127	ng/u1	98
30) Naphthalene	10.839	128	371758	4.612	ng/u1	99
33) Hexachlorobutadiene	11.127	225	55773	4.172	ng/u1	98
35) 4-Chloro-3-methylphenol	12.069	107	104371	4.326	ng/u1	98
36) 2-Methylnaphthalene	12.445	142	243250	4.529	ng/u1	99
37) 1-Methylnaphthalene	12.663	142	245545	4.548	ng/u1	100
39) 1,2,4,5-Tetrachloroben...	12.810	216	112652	4.086	ng/u1	98
41) 2,4,6-Trichlorophenol	13.045	196	65176	3.861	ng/u1	99
42) 2,4,5-Trichlorophenol	13.116	196	72646	4.014	ng/u1	99
43) 1,1'-Biphenyl	13.445	154	321312	4.475	ng/u1	100
44) 2-Chloronaphthalene	13.492	162	244407	4.338	ng/u1	98
45) 2-Nitroaniline	13.692	65	65634	4.733	ng/u1	97
47) Dimethylphthalate	14.068	163	313292	4.496	ng/u1	98
48) 2,6-Dinitrotoluene	14.186	165	44828	3.957	ng/u1	90
50) Acenaphthylene	14.333	152	367188	4.121	ng/u1	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092922\
 Data File : BM036796.D
 Acq On : 29 Sep 2022 11:47
 Operator : CG/JU
 Sample : SSTD00578
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD005026

Quant Time: Sep 30 01:11:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 30 01:10:23 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Acenaphthene	14.674	153	284854	4.569	ng/ul	99
56) Dibenzofuran	15.004	168	383762	4.584	ng/ul	99
57) 2,4-Dinitrotoluene	14.963	165	67058	4.071	ng/ul	90
58) 2,3,4,6-Tetrachlorophenol	15.227	232	58984	3.893	ng/ul#	99
59) Diethylphthalate	15.421	149	324286	4.648	ng/ul	99
61) Fluorene	15.651	166	317304	4.450	ng/ul	97
62) 4-Chlorophenyl-phenyle...	15.645	204	148492	4.355	ng/ul	97
67) N-Nitrosodiphenylamine	15.857	169	264068	4.481	ng/ul	99
68) 4-Bromophenyl-phenylether	16.539	248	79695	4.041	ng/ul	98
69) Hexachlorobenzene	16.645	284	94013	4.030	ng/ul	97
72) Phenanthrene	17.386	178	513356	4.658	ng/ul	99
74) Anthracene	17.480	178	498961	4.454	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.410	216	114180	4.172	ng/uL	99
76) Pentachlorobenzene	14.921	250	123760	4.181	ng/uL	97
78) Di-n-butylphthalate	18.309	149	504505	4.291	ng/ul	100
80) Fluoranthene	19.392	202	529244	4.415	ng/ul	99
82) Pyrene	19.756	202	585995	4.618	ng/ul	99
83) Butylbenzylphthalate	20.650	149	198554	4.130	ng/ul	93
85) Benzo(a)anthracene	21.497	228	567453	4.555	ng/ul	100
86) Bis(2-ethylhexyl)phtha...	21.427	149	312973	4.456	ng/ul	99
87) Chrysene	21.550	228	540166	4.568	ng/ul	99
90) Benzo(b)fluoranthene	23.191	252	556372	4.407	ng/ul	99
91) Benzo(k)fluoranthene	23.238	252	545350	4.323	ng/ul	100
93) Benzo(a)pyrene	23.821	252	475146	4.282	ng/ul	98
94) Indeno(1,2,3-cd)pyrene	26.444	276	566816	4.374	ng/ul	99
95) Dibenzo(a,h)anthracene	26.456	278	515644	4.387	ng/ul	99
96) Benzo(g,h,i)perylene	27.209	276	499989	4.450	ng/ul	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092922\
Data File : BM036796.D
Acq On : 29 Sep 2022 11:47
Operator : CG/JU
Sample : SST00578
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTD005026

Quant Time: Sep 30 01:11:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092922.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Sep 30 01:10:23 2022
Response via : Initial Calibration

