

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022718\
 Data File : BM014381.D
 Acq On : 27 Feb 2018 14:12
 Operator : SJ/JU
 Sample : J1631-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5Y7

Quant Time: Feb 27 16:40:13 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 27 03:22:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	109492	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	436508	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.17	164	244404	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.94	188	500495	20.00	ng/ul	0.00
75) Chrysene-d12	21.16	240	506467	20.00	ng/ul	0.00
83) Perylene-d12	23.33	264	484825	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.10	96	9348	3.65	ng/uL	0.00
5) Phenol-d5	6.72	99	197426	21.33	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	124887	22.29	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	167742	23.64	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	149858	18.51	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	82565	25.60	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	96481	28.90	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	164914	23.85	ng/ul	0.00
29) 4-Chloroaniline-d4	10.46	131	94587	13.88	ng/ul	0.00
43) Dimethylphthalate-d6	13.60	166	524618	25.99	ng/ul	0.00
46) Acenaphthylene-d8	13.86	160	668351	26.85	ng/ul	0.00
51) 4-Nitrophenol-d4	14.45	143	51886	18.62	ng/ul	0.02
57) Fluorene-d10	15.18	176	457653	25.31	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.34	200	43684	14.55	ng/ul	0.00
70) Anthracene-d10	17.04	188	677262	28.05	ng/ul	0.00
76) Pyrene-d10	19.35	212	721007	27.60	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.20	264	640667	27.15	ng/ul	0.02

Target Compounds

					Ovalue	
6) Phenol	6.75	94	21236	2.187	ng/ul	93
14) Acetophenone	8.35	105	36788	2.633	ng/ul#	96
28) Naphthalene	10.33	128	146800	6.548	ng/ul	97
34) 2-Methylnaphthalene	11.96	142	45572	2.671	ng/ul	95
44) Dimethylphthalate	13.65	163	215968	10.851	ng/ul#	97
47) Acenaphthylene	13.90	152	146639	6.234	ng/ul	97
49) Acenaphthene	14.24	153	112904	6.816	ng/ul	95
53) Dibenzofuran	14.58	168	124177	5.000	ng/ul	95
58) Fluorene	15.23	166	149084	6.953	ng/ul	97
69) Phenanthrene	16.99	178	2782062	96.477	ng/ul	99
71) Anthracene	17.07	178	524034	18.070	ng/ul	98
72) Carbazole	17.36	167	251321	10.306	ng/ul	96
74) Fluoranthene	19.02	202	4600221	133.526	ng/ul#	92
77) Pyrene	19.38	202	4166921	123.800	ng/ul#	91
80) Benzo(a)anthracene	21.14	228	2678194	84.730	ng/ul	99
82) Chrysene	21.19	228	2419603	82.057	ng/ul	97
85) Benzo(b)fluoranthene	22.69	252	3095068	95.341	ng/ul#	98
86) Benzo(k)fluoranthene	22.73	252	1092736	35.404	ng/ul#	96
88) Benzo(a)pyrene	23.24	252	2224663	76.689	ng/ul#	98
89) Indeno(1,2,3-cd)pyrene	25.51	276	1315270	46.724	ng/ul#	96
90) Dibenzo(a,h)anthracene	25.50	278	360810	15.426	ng/ul#	89
91) Benzo(a,h,i)perylene	26.17	276	922897	40.471	ng/ul#	95

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

