

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
 Data File : BM014413.D
 Acq On : 28 Feb 2018 21:56
 Operator : SJ/JU
 Sample : J1678-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5T5

Manual Integrations
 APPROVED

Sohil
 3/1/2018 5:15:05 PM

Quant Time: Mar 01 01:49:42 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 01 00:47:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	74079	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	359286	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.17	164	245088	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.94	188	602514	20.00	ng/ul	0.00
75) Chrysene-d12	21.15	240	702635	20.00	ng/ul	0.00
83) Perylene-d12	23.33	264	730494	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.09	96	8399m	4.85	ng/uL	-0.01
5) Phenol-d5	6.72	99	171070	27.32	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	109444	28.88	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	144427	30.08	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	146202	26.69	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	78253	29.47	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	90095	32.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	151532	26.62	ng/ul	0.00
29) 4-Chloroaniline-d4	10.45	131	206541	36.83	ng/ul	0.00
43) Dimethylphthalate-d6	13.60	166	623546	30.81	ng/ul	0.00
46) Acenaphthylene-d8	13.86	160	770853	30.89	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	55081m	19.71	ng/ul	0.01
57) Fluorene-d10	15.18	176	533970	29.45	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.34	200	64485	17.84	ng/ul	0.00
70) Anthracene-d10	17.03	188	861001	29.62	ng/ul	0.00
76) Pyrene-d10	19.34	212	1040907	28.72	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.19	264	1075478	30.25	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.75	94	11105m	1.690	ng/ul
28) Naphthalene	10.33	128	27892	1.511	ng/ul# 92
44) Dimethylphthalate	13.64	163	114396	5.732	ng/ul 99
53) Dibenzofuran	14.59	168	44864	1.801	ng/ul 90
58) Fluorene	15.24	166	48432	2.252	ng/ul 97
69) Phenanthrene	16.98	178	832905	23.993	ng/ul 98
71) Anthracene	17.07	178	154215	4.417	ng/ul 96
72) Carbazole	17.36	167	55198	1.880	ng/ul 96
74) Fluoranthene	19.01	202	979244	23.611	ng/ul# 94
77) Pyrene	19.37	202	820643	17.574	ng/ul# 93
80) Benzo(a)anthracene	21.13	228	387229	8.830	ng/ul 96
82) Chrysene	21.19	228	335399	8.199	ng/ul 93
85) Benzo(b)fluoranthene	22.68	252	438288	8.961	ng/ul# 97
86) Benzo(k)fluoranthene	22.71	252	123270m	2.651	ng/ul
88) Benzo(a)pyrene	23.23	252	312357	7.146	ng/ul# 99
89) Indeno(1,2,3-cd)pyrene	25.49	276	211496	4.987	ng/ul 98
90) Dibenzo(a,h)anthracene	25.49	278	53992m	1.532	ng/ul
91) Benzo(g,h,i)perylene	26.14	276	161713	4.707	ng/ul# 92

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

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