

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062916\
 Data File : BM006132.D
 Acq On : 29 Jun 2016 19:15
 Operator : UM/SJ
 Sample : H3779-20
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9YE9

Quant Time: Jun 30 03:43:11 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 29 00:57:09 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	208141	20.00	ng/ul	0.00
7) Naphthalene-d8	10.43	136	1015707	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	635057	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	1498051	20.00	ng/ul	0.00
29) Chrysene-d12	21.24	240	1798882	20.00	ng/ul	0.00
34) Perylene-d12	23.46	264	1696566	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.17	96	5351	1.19	ng/uL	0.00
3) Phenol-d5	6.82	99	272998	16.05	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	174489	17.40	ng/ul	0.00
5) 2-Chlorophenol-d4	7.18	132	207829	15.92	ng/ul	0.00
6) 4-Methylphenol-d8	8.35	113	217617	15.90	ng/ul	-0.01
8) Nitrobenzene-d5	8.80	128	111286	16.27	ng/ul	0.00
9) 2-Nitrophenol-d4	9.52	143	119627	15.28	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	224827	15.74	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	275032	24.36	ng/ul	0.00
16) Dimethylphthalate-d6	13.71	166	759357	16.68	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	940435	17.02	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	120711	13.98	ng/ul	0.00
21) Fluorene-d10	15.30	176	670412	17.26	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	95136	17.93	ng/ul	0.00
26) Anthracene-d10	17.14	188	1048201	17.44	ng/ul	0.00
30) Pyrene-d10	19.44	212	1228837	16.96	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.31	264	1142321	16.92	ng/ul	0.00

Target Compounds

						Qvalue
28) Fluoranthene	19.11	202	192338	2.22	ng/ul	99
31) Pyrene	19.47	202	162508	1.71	ng/ul	98
32) Benzo(a)anthracene	21.23	228	98257	1.05	ng/ul	97
33) Chrysene	21.28	228	119940	1.39	ng/ul	98
35) Benzo(b)fluoranthene	22.79	252	172152	1.96	ng/ul	98
38) Benzo(a)pyrene	23.36	252	100767	1.21	ng/ul	98

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

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