

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062616\
 Data File : BM006039.D
 Acq On : 26 Jun 2016 14:24
 Operator : UM/SJ
 Sample : H3669-19
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y72

Quant Time: Jun 27 05:33:17 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 27 05:15:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	259568	20.00	ng/ul	0.00
7) Naphthalene-d8	10.43	136	1302609	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	854131	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	2107067	20.00	ng/ul	0.00
29) Chrysene-d12	21.25	240	2365614	20.00	ng/ul	0.00
34) Perylene-d12	23.47	264	2259391	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.18	96	4530	0.78	ng/uL	0.00
3) Phenol-d5	6.83	99	257927	9.45	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	160480	10.56	ng/ul	0.00
5) 2-Chlorophenol-d4	7.19	132	196159	10.31	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	180505	7.83	ng/ul	0.00
8) Nitrobenzene-d5	8.81	128	104019	10.49	ng/ul	0.00
9) 2-Nitrophenol-d4	9.52	143	112692	10.21	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	225329	10.64	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	180952	7.88	ng/ul	0.00
16) Dimethylphthalate-d6	13.72	166	794780	10.87	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	982119	11.58	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	104026	6.67	ng/ul	0.00
21) Fluorene-d10	15.30	176	736668	11.80	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	83219	7.11	ng/ul	0.00
26) Anthracene-d10	17.15	188	1176095	12.15	ng/ul	0.00
30) Pyrene-d10	19.45	212	1393681	13.45	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.33	264	1235626	12.11	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.09	178	204593	1.81	ng/ul	99
28) Fluoranthene	19.12	202	378797	2.61	ng/ul	99
31) Pyrene	19.48	202	321718	2.38	ng/ul	100
32) Benzo(a)anthracene	21.23	228	197281	1.40	ng/ul	94
33) Chrysene	21.29	228	250401	1.95	ng/ul#	95
35) Benzo(b)fluoranthene	22.80	252	357563	2.72	ng/ul#	89
38) Benzo(a)pyrene	23.37	252	202923	1.61	ng/ul#	93
39) Indeno(1,2,3-cd)pyrene	25.69	276	151503	1.01	ng/ul	98
41) Benzo(g,h,i)perylene	26.37	276	146480	1.13	ng/ul	94

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

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