

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062816\
 Data File : BM006105.D
 Acq On : 28 Jun 2016 20:50
 Operator : UM/SJ
 Sample : PB91650BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK50

Quant Time: Jun 29 00:59:15 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	179245	20.00	ng/ul	0.00
7) Naphthalene-d8	10.43	136	781706	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	454620	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	953366	20.00	ng/ul	0.00
29) Chrysene-d12	21.24	240	923083	20.00	ng/ul	0.00
34) Perylene-d12	23.46	264	919302	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	3.16	96	14597	3.78	ng/uL	0.00
3) Phenol-d5	6.83	99	600935	41.04	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	381179	44.15	ng/ul	0.00
5) 2-Chlorophenol-d4	7.19	132	464407	41.32	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	481703	40.87	ng/ul	0.00
8) Nitrobenzene-d5	8.80	128	239132	45.42	ng/ul	0.00
9) 2-Nitrophenol-d4	9.52	143	263428	43.71	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	448073	40.75	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	662767	76.26	ng/ul	0.00
16) Dimethylphthalate-d6	13.72	166	1441381	44.23	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	1801398	45.55	ng/ul	0.00
20) 4-Nitrophenol-d4	14.51	143	241719	39.09	ng/ul	0.00
21) Fluorene-d10	15.30	176	1205781	43.36	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	157376	46.61	ng/ul	0.00
26) Anthracene-d10	17.15	188	1782075	46.60	ng/ul	0.00
30) Pyrene-d10	19.45	212	1833343	49.32	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.32	264	1728572	47.25	ng/ul	0.00

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

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