

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062416\
 Data File : BM005969.D
 Acq On : 24 Jun 2016 17:13
 Operator : UM/SJ
 Sample : PB91577TB
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB17

Quant Time: Jun 25 00:29:07 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 23:55:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	169436	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	735083	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.31	164	408273	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	919747	20.00	ng/ul	0.00
75) Chrysene-d12	21.24	240	1094563	20.00	ng/ul	0.00
83) Perylene-d12	23.46	264	1128518	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.18	96	21279	5.00	ng/uL	0.00
5) Phenol-d5	6.83	99	699290	47.92	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	443191	49.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	549665	47.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	539908	47.23	ng/ul	0.00
19) Nitrobenzene-d5	8.81	128	255192	47.19	ng/ul	0.00
22) 2-Nitrophenol-d4	9.53	143	273516	48.32	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	496201	45.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	733950	61.86	ng/ul	0.00
43) Dimethylphthalate-d6	13.72	166	1506890	46.45	ng/ul	0.00
46) Acenaphthylene-d8	14.00	160	1915613	47.28	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	270675	42.68	ng/ul	0.00
57) Fluorene-d10	15.30	176	1305947	46.60	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.42	200	187871	41.01	ng/ul	0.00
70) Anthracene-d10	17.15	188	2020544	48.23	ng/ul	0.00
76) Pyrene-d10	19.45	212	2285860	45.00	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.33	264	2464449	48.47	ng/ul	0.00

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

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