

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110716\
 Data File : BM007773.D
 Acq On : 07 Nov 2016 16:12
 Operator : UM/SJ
 Sample : PB94579BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK79

Quant Time: Nov 07 17:32:10 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 24 12:55:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	194099	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	742623	20.00	ng/ul	-0.01
35) Acenaphthene-d10	14.38	164	470611	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.13	188	907760	20.00	ng/ul	0.00
75) Chrysene-d12	21.32	240	1101305	20.00	ng/ul	0.00
83) Perylene-d12	23.57	264	1124637	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	30929	6.83	ng/uL	0.00
5) Phenol-d5	6.92	99	538209	35.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.09	67	322129	36.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.28	132	427783	37.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	425035	36.26	ng/ul	0.00
19) Nitrobenzene-d5	8.90	128	203206	38.02	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	222827	39.33	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	401057	36.01	ng/ul	0.00
29) 4-Chloroaniline-d4	10.67	131	527910	41.55	ng/ul	0.00
43) Dimethylphthalate-d6	13.79	166	1220537	38.09	ng/ul	0.00
46) Acenaphthylene-d8	14.07	160	1448343	37.55	ng/ul	0.00
51) 4-Nitrophenol-d4	14.60	143	158563	31.30	ng/ul	0.00
57) Fluorene-d10	15.37	176	1027110	36.57	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.51	200	155500	28.83	ng/ul	0.00
70) Anthracene-d10	17.23	188	1571164	38.41	ng/ul	0.00
76) Pyrene-d10	19.52	212	1782389	39.39	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.43	264	1899042	38.68	ng/ul	0.00

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

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