

Data Path : Z:\HPCHEM1\BNA M\DATA\BM091916\
 Data File : BM007479.D
 Acq On : 19 Sep 2016 18:02
 Operator : UM/SJ
 Sample : PB93450BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK50

Quant Time: Sep 20 01:44:00 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 20 01:35:12 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	161725	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	745868	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.42	164	586746	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1333343	20.00	ng/ul	0.00
75) Chrysene-d12	21.34	240	1888523	20.00	ng/ul	0.00
83) Perylene-d12	23.62	264	1927040	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.30	96	26637	8.42	ng/uL	0.00
5) Phenol-d5	6.96	99	528968	34.62	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	326627	40.69	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	413966	36.66	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	458981	34.45	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	219007	40.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	258260	40.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	447603	35.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.71	131	495744	31.22	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	1802843	35.62	ng/ul	0.00
46) Acenaphthylene-d8	14.11	160	2054494	35.02	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	286840	30.02	ng/ul	0.00
57) Fluorene-d10	15.42	176	1501606	34.30	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	284737	33.96	ng/ul	0.00
70) Anthracene-d10	17.26	188	2545554	40.87	ng/ul	0.00
76) Pyrene-d10	19.56	212	3166722	40.10	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.47	264	3554249	40.04	ng/ul	0.00

Target Compounds	Ovalue

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

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