

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092316\
 Data File : BM007547.D
 Acq On : 24 Sep 2016 05:25
 Operator : UM/SJ
 Sample : PB93566BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK66

Quant Time: Sep 24 06:25:46 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 24 04:39:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	222377	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	1019947	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	816114	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1882822	20.00	ng/ul	0.00
75) Chrysene-d12	21.34	240	2616323	20.00	ng/ul	0.00
83) Perylene-d12	23.60	264	2539809	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	35486	7.84	ng/uL	0.00
5) Phenol-d5	6.95	99	683510	34.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	411193	38.75	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	536595	37.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	599851	35.60	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	298117	40.88	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	351805	42.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	606741	36.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	855544	49.37	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	2482880	41.83	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	2744956	39.85	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	408106	38.95	ng/ul	0.00
57) Fluorene-d10	15.41	176	2068253	39.73	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	420220	36.58	ng/ul	0.00
70) Anthracene-d10	17.26	188	3484053	40.27	ng/ul	0.00
76) Pyrene-d10	19.55	212	4355628	41.95	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.46	264	4631582	41.21	ng/ul	0.00

Target Compounds	Ovalue

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

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