

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111416\
 Data File : BM007849.D
 Acq On : 14 Nov 2016 11:32
 Operator : UM/SJ
 Sample : H5610-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BD4M8

Quant Time: Nov 15 02:27:14 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 15 02:14:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	151817	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	581301	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	368545	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	735439	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	910450	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	935562	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.28	96	25210	7.11	ng/uL	0.00
5) Phenol-d5	6.91	99	469345	40.06	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	286411	41.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	373723	41.59	ng/ul	0.00
13) 4-Methylphenol-d8	8.44	113	371212	40.49	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	181824	43.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	196860	44.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	352427	40.42	ng/ul	0.00
29) 4-Chloroaniline-d4	10.66	131	361420	36.34	ng/ul	0.00
43) Dimethylphthalate-d6	13.79	166	1092483	43.54	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1293555	42.83	ng/ul	0.00
51) 4-Nitrophenol-d4	14.59	143	143004	36.04	ng/ul	0.00
57) Fluorene-d10	15.37	176	945216	42.98	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.50	200	132480	30.32	ng/ul	0.00
70) Anthracene-d10	17.22	188	1463396	44.16	ng/ul	0.00
76) Pyrene-d10	19.52	212	1696145	45.34	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.41	264	1840294	45.06	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Phenol	6.94	94	55294	4.60	ng/ul	96
44) Dimethylphthalate	13.83	163	171301	7.03	ng/ul	100

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

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