

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102016\
 Data File : BM007710.D
 Acq On : 20 Oct 2016 19:15
 Operator : UM/SJ
 Sample : H5330-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B84

Quant Time: Oct 21 10:13:16 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 17:01:48 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	163267	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.53	136	610523	20.00	ng/ul	-0.02
35) Acenaphthene-d10	14.39	164	382795	20.00	ng/ul	-0.02
61) Phenanthrene-d10	17.13	188	759677	20.00	ng/ul	-0.02
75) Chrysene-d12	21.32	240	963933	20.00	ng/ul	-0.02
83) Perylene-d12	23.57	264	994757	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	5893	1.55	ng/uL	0.00
5) Phenol-d5	6.93	99	86421	6.86	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	7.09	67	257434	34.43	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.29	132	274827	28.44	ng/ul	-0.02
13) 4-Methylphenol-d8	8.46	113	164230	16.66	ng/ul	-0.02
19) Nitrobenzene-d5	8.91	128	159557	36.32	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.63	143	172244	36.98	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.16	165	300341	32.80	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.70	131	10832	1.04	ng/ul	0.00
43) Dimethylphthalate-d6	13.80	166	995611	38.20	ng/ul	-0.02
46) Acenaphthylene-d8	14.08	160	1186152	37.81	ng/ul	-0.02
51) 4-Nitrophenol-d4	14.62	143	16759	4.07	ng/ul	0.00
57) Fluorene-d10	15.38	176	864468	37.84	ng/ul	-0.02
62) 4,6-Dinitro-2-methylphenol	15.51	200	133775	29.64	ng/ul	-0.02
70) Anthracene-d10	17.23	188	1328562	38.81	ng/ul	-0.02
76) Pyrene-d10	19.53	212	1531580	38.67	ng/ul	-0.02
87) Benzo(a)pyrene-d12	23.43	264	1680449	38.69	ng/ul	-0.02

Target Compounds	Ovalue

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

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