

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121816\
 Data File : BM008426.D
 Acq On : 17 Dec 2016 23:50
 Operator : UM/SJ
 Sample : H6039-11DL 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8S7DL

Quant Time: Dec 19 16:51:50 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Dec 18 04:15:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	110741	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	436980	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.30	164	292854	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	591160	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	735457	20.00	ng/ul	0.00
83) Perylene-d12	23.49	264	739204	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.89	99	11085	1.29	ng/ul	0.04
7) Bis-(2-Chloroethyl)ether-d	7.01	67	29909	6.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.21	132	34468	5.33	ng/ul	0.01
13) 4-Methylphenol-d8	8.39	113	23846	3.58	ng/ul	0.02
19) Nitrobenzene-d5	8.83	128	20449	6.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	23870	6.82	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	43698	6.81	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	13.72	166	151847	7.80	ng/ul	0.00
46) Acenaphthylene-d8	13.99	160	174544	7.59	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.30	176	128246	7.58	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.48	200	14044	3.99	ng/ul	0.02
70) Anthracene-d10	17.16	188	202849	7.88	ng/ul	0.00
76) Pyrene-d10	19.46	212	241512	8.50	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.34	264	241830	7.79	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
24) 2,4-Dimethylphenol	9.63	107	244278	30.56	ng/ul	94
28) Naphthalene	10.48	128	39755	1.92	ng/ul	98

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

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