

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041516\
 Data File : BM004958.D
 Acq On : 15 Apr 2016 14:18
 Operator : UM/SJ
 Sample : H2298-16
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9T98

Quant Time: Apr 15 23:18:06 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM041416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 15 22:59:24 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	55711	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	255081	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.45	164	162256	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	377388	20.00	ng/ul	0.00
75) Chrysene-d12	21.38	240	391734	20.00	ng/ul	0.00
83) Perylene-d12	23.67	264	318718	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.31	96	1232	1.18	ng/uL	0.00
5) Phenol-d5	6.97	99	23661	5.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	78099	34.24	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	82498	24.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	50471	15.36	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	55485	32.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.69	143	62372	32.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	100633	29.14	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	13.85	166	397226	35.99	ng/ul	0.00
46) Acenaphthylene-d8	14.14	160	463130	34.21	ng/ul	0.00
51) 4-Nitrophenol-d4	14.65	143	5879	3.12	ng/ul	0.00
57) Fluorene-d10	15.44	176	338458	35.50	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.56	200	45938	28.14	ng/ul	0.00
70) Anthracene-d10	17.29	188	523783	34.93	ng/ul	0.00
76) Pyrene-d10	19.59	212	581042	35.59	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.52	264	446948	36.41	ng/ul	0.00

Target Compounds Ovalue

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

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