

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041216\
 Data File : BM004943.D
 Acq On : 12 Apr 2016 15:59
 Operator : UM/SJ
 Sample : H1935-12
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AH0

Quant Time: Apr 12 18:43:23 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM040116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 12 18:19:19 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	38831	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	179076	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.45	164	115593	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	277893	20.00	ng/ul	0.00
78) Chrysene-d12	21.38	240	343820	20.00	ng/ul	0.00
86) Perylene-d12	23.67	264	331357	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.31	96	6339	6.60	ng/uL	0.00
5) Phenol-d5	6.98	99	124043	35.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	80027	42.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	102160	39.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	104572	36.09	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	51430	40.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.69	143	57773	40.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	102530	38.93	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	2106	0.67	ng/ul	0.01
44) Dimethylphthalate-d6	13.86	166	406023	44.52	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	447795	41.05	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	66296	36.77	ng/ul	0.00
58) Fluorene-d10	15.45	176	344275	43.68	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.56	200	61906	37.18	ng/ul	0.00
71) Anthracene-d10	17.30	188	564011	45.36	ng/ul	0.00
79) Pyrene-d10	19.59	212	673957	45.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.52	264	755546	51.47	ng/ul	0.00

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

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