

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040616\
 Data File : BM004902.D
 Acq On : 06 Apr 2016 13:24
 Operator : UM/SJ
 Sample : PB89480BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK80

Quant Time: Apr 07 04:32:21 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM040116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 04:29:45 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	57582	20.00	ng/ul	0.00
18) Naphthalene-d8	10.61	136	261286	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	164768	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	406828	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	526781	20.00	ng/ul	0.00
86) Perylene-d12	23.68	264	538354	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.31	96	6194	4.35	ng/uL	0.00
5) Phenol-d5	6.98	99	109723	21.44	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	68384	24.66	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	90029	23.18	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	90877	21.15	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	42891	23.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	48232	23.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	85213	22.17	ng/ul	0.00
29) 4-Chloroaniline-d4	10.75	131	133439	29.09	ng/ul	0.00
44) Dimethylphthalate-d6	13.86	166	336234	25.87	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	379661	24.41	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	51546	20.06	ng/ul	0.00
58) Fluorene-d10	15.45	176	287356	25.58	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	45411	18.63	ng/ul	0.00
71) Anthracene-d10	17.30	188	469442	25.79	ng/ul	0.00
79) Pyrene-d10	19.59	212	570121	24.93	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.53	264	635921	26.66	ng/ul	0.00

Target Compounds Ovalue

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

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