

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100115\
 Data File : BM002805.D
 Acq On : 01 Oct 2015 15:45
 Operator : UM/IZ
 Sample : PB85824BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 02 01:22:18 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 02 01:21:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.69	152	66060	20.00	ng/ul	0.00
18) Naphthalene-d8	11.55	136	272700	20.00	ng/ul	0.00
36) Acenaphthene-d10	15.29	164	136829	20.00	ng/ul	0.00
62) Phenanthrene-d10	18.00	188	278918	20.00	ng/ul	0.00
78) Chrysene-d12	22.09	240	299397	20.00	ng/ul	0.00
86) Perylene-d12	24.67	264	304875	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.76	96	9247	6.50	ng/uL	0.00
5) Phenol-d5	7.80	99	166427	33.25	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.98	67	93379	33.16	ng/ul	0.00
9) 2-Chlorophenol-d4	8.20	132	154072	33.07	ng/ul	0.00
13) 4-Methylphenol-d8	9.39	113	137453	33.31	ng/ul	0.00
19) Nitrobenzene-d5	9.88	128	71321	34.30	ng/ul	0.00
22) 2-Nitrophenol-d4	10.62	143	75875	34.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	11.16	165	124168	32.37	ng/ul	0.00
29) 4-Chloroaniline-d4	11.67	131	196210	42.49	ng/ul	0.00
44) Dimethylphthalate-d6	14.66	166	370530	35.99	ng/ul	0.00
47) Acenaphthylene-d8	14.99	160	475397	34.60	ng/ul	0.00
52) 4-Nitrophenol-d4	15.46	143	62589	34.24	ng/ul	0.00
58) Fluorene-d10	16.26	176	312513	33.70	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	16.38	200	40111	29.49	ng/ul	0.00
71) Anthracene-d10	18.09	188	465436	34.97	ng/ul	0.00
79) Pyrene-d10	20.33	212	488714	34.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.50	264	562022	36.04	ng/ul	0.00

Target Compounds Ovalue

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

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