

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121115\
 Data File : BM003494.D
 Acq On : 11 Dec 2015 22:31
 Operator : UM/IZ
 Sample : PB87221BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK21

Quant Time: Dec 12 01:27:16 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM120115.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 12 00:54:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	1597	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	8027	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	4587	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	7976	20.00	ng/ul	0.00
78) Chrysene-d12	21.32	240	7650	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	7781	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.18	96	10653	331.46	ng/uL	0.00
5) Phenol-d5	6.89	99	196489	1541.13	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	112744	1602.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	168508	1598.54	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	163598	1529.54	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	81028	1479.90	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	86668	1514.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	152693	1345.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	234424	1825.40	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	452730	1312.82	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	569606	1343.77	ng/ul	0.00
52) 4-Nitrophenol-d4	14.57	143	56493	913.94	ng/ul	0.00
58) Fluorene-d10	15.36	176	391032	1317.72	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	39080	957.67	ng/ul	0.00
71) Anthracene-d10	17.22	188	549829	1551.61	ng/ul	0.00
79) Pyrene-d10	19.52	212	568626	1518.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.43	264	597556	1591.72	ng/ul	0.00

Target Compounds	Ovalue

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

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