

Data Path : Z:\HPCHEM1\BNA G\DATA\BG090115\
 Data File : BG018559.D
 Acq On : 1 Sep 2015 12:28
 Operator : UM/IZ
 Sample : PB85282BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK82

Manual Integrations
 APPROVED

MMDadoda
 9/2/2015 2:06:01 PM

Quant Time: Sep 02 02:32:59 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 02 02:18:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	78092	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	344233	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.64	164	239543	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.38	188	643670	20.00	ng/ul	0.00
78) Chrysene-d12	21.64	240	688291	20.00	ng/ul	0.00
86) Perylene-d12	24.84	264	671908	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.45	96	11793	6.67	ng/uL	0.00
5) Phenol-d5	7.17	99	232077	32.74	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	143289	32.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	174246	32.10	ng/ul	0.00
13) 4-Methylphenol-d8	8.71	113	193739	32.53	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	94000	35.03	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	102161	37.38	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.43	165	197761	34.10	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	287840	43.91	ng/ul	0.00
44) Dimethylphthalate-d6	14.04	166	768116	38.11	ng/ul	0.00
47) Acenaphthylene-d8	14.33	160	819943	32.31	ng/ul	0.00
52) 4-Nitrophenol-d4	14.82	143	131611m	36.30	ng/ul	0.00
58) Fluorene-d10	15.63	176	607494	34.61	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	82613	26.09	ng/ul	-0.01
71) Anthracene-d10	17.47	188	1031402	33.50	ng/ul	0.00
79) Pyrene-d10	19.76	212	1139729	31.92	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.62	264	1231627	34.53	ng/ul	0.00

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

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