

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080315\
 Data File : BG018243.D
 Acq On : 3 Aug 2015 14:49
 Operator : UM/TP
 Sample : PB84804BL
 Misc : SBLK04
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK04

Manual Integrations
 APPROVED

MMDadoda
 8/4/2015 12:16:55 PM

Quant Time: Aug 04 03:01:41 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG073115.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 04 02:54:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	32085	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	145750	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	99896	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	245340m	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	238093m	20.00	ng/ul	0.00
86) Perylene-d12	23.27	264	164168	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.94	96	3297	7.76	ng/uL	0.00
5) Phenol-d5	6.66	99	90155	31.25	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.80	67	47045	31.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	72983	33.49	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	75391	32.09	ng/ul	0.00
19) Nitrobenzene-d5	8.61	128	40929	35.65	ng/ul	0.00
22) 2-Nitrophenol-d4	9.33	143	47315	35.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.88	165	76816	32.99	ng/ul	0.00
29) 4-Chloroaniline-d4	10.39	131	115615	39.84	ng/ul	0.00
44) Dimethylphthalate-d6	13.55	166	302832	39.23	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	321838	35.96	ng/ul	0.00
52) 4-Nitrophenol-d4	14.39	143	46012	32.20	ng/ul	0.00
58) Fluorene-d10	15.14	176	249819	35.76	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.28	200	45925m	26.81	ng/ul	0.00
71) Anthracene-d10	17.00	188	411750	38.60	ng/ul	0.00
79) Pyrene-d10	19.31	212	437949	38.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.14	264	330098	39.76	ng/ul	0.00

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

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