

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
 Data File : BG024229.D
 Acq On : 7 Oct 2016 11:28
 Operator : UM/SJ
 Sample : PB93836BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK36

Manual Integrations
 APPROVED

sohil
 10/7/2016 7:00:59 PM

Quant Time: Oct 07 13:37:35 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 07 06:34:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	136272	20.00	ng/ul	0.00
18) Naphthalene-d8	11.13	136	580600	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.91	164	483515	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.65	188	1034445m	20.00	ng/ul	0.00
75) Chrysene-d12	21.95	240	1098390	20.00	ng/ul	0.00
83) Perylene-d12	25.41	264	1036547	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.65	96	17891m	6.23	ng/uL	-0.01
5) Phenol-d5	7.42	99	373506	29.38	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.61	67	266128m	30.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.82	132	265772	30.36	ng/ul	0.00
13) 4-Methylphenol-d8	8.99	113	308349	30.00	ng/ul	0.00
19) Nitrobenzene-d5	9.47	128	136424	33.11	ng/ul	-0.01
22) 2-Nitrophenol-d4	10.20	143	155942	33.83	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.73	165	299386	29.82	ng/ul	0.00
29) 4-Chloroaniline-d4	11.25	131	413144	39.17	ng/ul	0.00
43) Dimethylphthalate-d6	14.31	166	1123095	33.27	ng/ul	0.00
46) Acenaphthylene-d8	14.61	160	1263166	32.66	ng/ul	0.00
51) 4-Nitrophenol-d4	15.08	143	182848	30.54	ng/ul	0.00
57) Fluorene-d10	15.90	176	1033631	32.38	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.00	200	188024m	30.24	ng/ul	0.00
70) Anthracene-d10	17.75	188	1585702m	33.94	ng/ul	0.00
76) Pyrene-d10	20.02	212	1750483	34.14	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.17	264	1619107	34.72	ng/ul	0.00

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

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