

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101216\
 Data File : BG024310.D
 Acq On : 12 Oct 2016 20:45
 Operator : UM/SJ
 Sample : PB93938BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK38

Manual Integrations
 APPROVED

sohil
 10/13/2016 1:55:14 PM

Quant Time: Oct 13 01:53:09 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 13 01:34:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	134198	20.00	ng/ul	0.00
18) Naphthalene-d8	11.12	136	597620	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.90	164	513405	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.65	188	1055003	20.00	ng/ul	0.00
75) Chrysene-d12	21.94	240	1163702	20.00	ng/ul	0.00
83) Perylene-d12	25.38	264	1192399	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.64	96	19558m	6.92	ng/uL	0.00
5) Phenol-d5	7.42	99	356933	28.51	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	233305	27.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	248916	28.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	311002	30.73	ng/ul	0.00
19) Nitrobenzene-d5	9.47	128	137391	32.39	ng/ul	0.00
22) 2-Nitrophenol-d4	10.19	143	160405	33.81	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	299368	28.97	ng/ul	0.00
29) 4-Chloroaniline-d4	11.24	131	402837	37.10	ng/ul	0.00
43) Dimethylphthalate-d6	14.30	166	1103264	30.78	ng/ul	0.00
46) Acenaphthylene-d8	14.60	160	1251687	30.48	ng/ul	0.00
51) 4-Nitrophenol-d4	15.07	143	178088	28.02	ng/ul	0.00
57) Fluorene-d10	15.89	176	1040891	30.71	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	193495	30.52	ng/ul	0.00
70) Anthracene-d10	17.75	188	1535147m	32.21	ng/ul	0.00
76) Pyrene-d10	20.01	212	1729984	31.85	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.15	264	1773126	33.05	ng/ul	0.00

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

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