

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111616\
 Data File : BG024797.D
 Acq On : 16 Nov 2016 10:36
 Operator : UM/SJ
 Sample : PB94741BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK41

Manual Integrations
 APPROVED

sohil
 11/17/2016 5:39:27 PM

Quant Time: Nov 17 02:24:42 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 17 01:26:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	87679	20.00	ng/ul	0.00
18) Naphthalene-d8	11.09	136	366196	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.88	164	298166	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.62	188	733949m	20.00	ng/ul	0.00
75) Chrysene-d12	21.92	240	1048226	20.00	ng/ul	0.00
83) Perylene-d12	25.34	264	1079811	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.62	96	13275	7.85	ng/uL	0.00
5) Phenol-d5	7.39	99	272377	36.26	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.57	67	176863	37.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.79	132	202911	37.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.95	113	220152	34.90	ng/ul	0.00
19) Nitrobenzene-d5	9.44	128	100225	36.15	ng/ul	0.00
22) 2-Nitrophenol-d4	10.16	143	113900	34.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.70	165	215116	32.65	ng/ul	0.00
29) 4-Chloroaniline-d4	11.22	131	273560	37.17	ng/ul	0.00
43) Dimethylphthalate-d6	14.27	166	813538	37.42	ng/ul	0.00
46) Acenaphthylene-d8	14.57	160	864492	36.49	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	140848	36.41	ng/ul	0.00
57) Fluorene-d10	15.87	176	736400	35.99	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.97	200	164509m	32.60	ng/ul	0.00
70) Anthracene-d10	17.72	188	1244638	38.39	ng/ul	0.00
76) Pyrene-d10	19.99	212	1641092	36.11	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.11	264	1849780	39.13	ng/ul	0.00

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

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