

Data Path : Z:\HPCHEM1\BNA_B\Data\BB101116\
 Data File : BB058603.D
 Acq On : 11 Oct 2016 11:41
 Operator : UM/SJ
 Sample : MDL-BLK-S
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_B
 ClientSampleId :
 MDL-BLK-S

Manual Integrations
 APPROVED

umangi
 10/11/2016 7:11:08 PM

Quant Time: Oct 11 14:12:19 2016
 Quant Method : Z:\HPCHEM1\BNA_B\METHOD\SOM02.2-EPA-BB101016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 14:09:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.77	152	560758	20.00	ng/ul	0.10
18) Naphthalene-d8	11.72	136	2077344	20.00	ng/ul	0.10
35) Acenaphthene-d10	15.65	164	1155352	20.00	ng/ul	0.06
61) Phenanthrene-d10	18.48	188	1781357m	20.00	ng/ul	0.04
75) Chrysene-d12	22.83	240	1766824	20.00	ng/ul	0.02
83) Perylene-d12	25.95	264	1468979	20.00	ng/ul	0.04

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.69	96	76127	6.15	ng/uL	0.06
5) Phenol-d5	7.91	99	1121160	27.52	ng/ul	0.10
7) Bis-(2-Chloroethyl)ether-d	8.06	67	816541	29.92	ng/ul	0.12
9) 2-Chlorophenol-d4	8.29	132	1182547	27.98	ng/ul	0.10
13) 4-Methylphenol-d8	9.54	113	898164	26.63	ng/ul	0.10
19) Nitrobenzene-d5	9.99	128	591910	27.56	ng/ul	0.10
22) 2-Nitrophenol-d4	10.76	143	647210	28.00	ng/ul	0.10
26) 2,4-Dichlorophenol-d3	11.33	165	868586	26.52	ng/ul	0.10
29) 4-Chloroaniline-d4	11.85	131	1257868	31.90	ng/ul	0.10
43) Dimethylphthalate-d6	15.03	166	2295611	29.88	ng/ul	0.06
46) Acenaphthylene-d8	15.32	160	2588648	28.57	ng/ul	0.06
51) 4-Nitrophenol-d4	15.82	143	383911	21.76	ng/ul	0.02
57) Fluorene-d10	16.67	176	1827197	29.03	ng/ul	0.06
62) 4,6-Dinitro-2-methylphenol	16.76	200	388268	20.53	ng/ul	0.02
70) Anthracene-d10	18.57	188	2309362	28.99	ng/ul	0.04
76) Pyrene-d10	20.90	212	2490837	32.08	ng/ul	0.02
87) Benzo(a)pyrene-d12	25.74	264	1919989	28.62	ng/ul	0.04

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

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