

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091718\
 Data File : BN002733.D
 Acq On : 17 Sep 2018 19:29
 Operator : JU/SJ
 Sample : J4893-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C08K3

Manual Integrations
 APPROVED

Sohil
 9/18/2018 6:16:47 PM

Quant Time: Sep 18 03:03:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 18 02:47:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	143437	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	566795	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.29	164	290465	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.03	188	590899	20.00	ng/ul	0.00
77) Chrysene-d12	21.24	240	595533	20.00	ng/ul	0.00
85) Perylene-d12	23.45	264	644329	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	14539	4.68	ng/uL	0.00
5) Phenol-d5	6.84	99	61657	5.06	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	164158	21.70	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	181223	18.08	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	105135	11.14	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	101997	23.23	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	113978	22.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	185086	20.64	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	21412	2.06	ng/ul	0.02
43) Dimethylphthalate-d6	13.70	166	551342	24.25	ng/ul	0.00
46) Acenaphthylene-d8	13.97	160	712028	23.98	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	13498m	3.59	ng/ul	0.04
57) Fluorene-d10	15.28	176	484321	24.31	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.43	200	69295	18.75	ng/ul	0.00
70) Anthracene-d10	17.13	188	711716	25.23	ng/ul	0.00
78) Pyrene-d10	19.43	212	746802	23.76	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.31	264	851508	25.11	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.87	94	26527	2.160	ng/ul 95
36) 1,2,4,5-Tetrachlorobenzene	12.46	216	139373	14.444	ng/ul 99
44) Dimethylphthalate	13.75	163	535148	24.116	ng/ul 100
72) 1,2,3,4-Tetrachlorobenzene	13.08	216	112439	11.138	ng/uL 99

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

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