

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM022983.D
 Acq On : 05 Oct 2019 10:08
 Operator : JU
 Sample : K5050-02
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLMBS

Quant Time: Oct 07 01:14:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 07 01:10:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	212819	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	931824	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	572593	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.16	188	1000890	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	735252	20.00	ng/ul	0.00
86) Perylene-d12	23.60	264	847693	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.27	96	11871	2.41	ng/uL	0.00
5) Phenol-d5	6.96	99	217640	11.57	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	126891	12.33	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	179037	11.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	179655	11.67	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	86463	12.02	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	87998	11.21	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	181274	11.59	ng/ul	0.00
29) 4-Chloroaniline-d4	10.71	131	240122	12.64	ng/ul	0.00
44) Dimethylphthalate-d6	13.83	166	611373	13.76	ng/ul	0.00
47) Acenaphthylene-d8	14.12	160	662559	12.91	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	49827	5.67	ng/ul	0.00
58) Fluorene-d10	15.42	176	490687	12.96	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	43158	6.63	ng/ul	0.00
71) Anthracene-d10	17.26	188	712776	14.58	ng/ul	0.00
79) Pyrene-d10	19.55	212	721881	18.24	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.46	264	707328	16.13	ng/ul	0.00

Target Compounds

				Ovalue
6) Phenol	6.98	94	38746	1.966 ng/ul 97
45) Dimethylphthalate	13.88	163	212718	4.660 ng/ul 99

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

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