

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091520\
 Data File : BM027417.D
 Acq On : 15 Sep 2020 18:16
 Operator : JU/CG
 Sample : L3995-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG0N7

Quant Time: Sep 15 18:49:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 15 13:17:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	120004	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	430289	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.20	164	279332	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.94	188	622201	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	707767	20.00	ng/ul	0.00
86) Perylene-d12	23.31	264	808615	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	17045	6.32	ng/uL	0.00
5) Phenol-d5	6.75	99	58626	5.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	141227	21.50	ng/ul	0.00
9) 2-Chlorophenol-d4	7.10	132	143032	19.75	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	100608	12.47	ng/ul	0.00
19) Nitrobenzene-d5	8.70	128	79412	24.52	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	86690	24.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	172731	22.59	ng/ul	0.00
29) 4-Chloroaniline-d4	10.46	131	197124	21.81	ng/ul	0.00
44) Dimethylphthalate-d6	13.62	166	612960	27.58	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	703899	28.04	ng/ul	0.00
52) 4-Nitrophenol-d4	14.43	143	7937	2.03	ng/ul	0.01
58) Fluorene-d10	15.20	176	524745	26.36	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	48271	12.12	ng/ul	0.00
71) Anthracene-d10	17.04	188	873937	29.70	ng/ul	0.00
79) Pyrene-d10	19.34	212	1074723	34.09	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.17	264	1256585	28.97	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.16	88	8460	2.621	ng/uL 98
6) Phenol	6.77	94	20533	1.953	ng/ul 92
45) Dimethylphthalate	13.66	163	94294	4.026	ng/ul 99

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

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