

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121820\
 Data File : BP004367.D
 Acq On : 18 Dec 2020 20:50
 Operator : CG/JU
 Sample : L5128-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8C0

Quant Time: Dec 19 01:01:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 18 17:12:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	540424	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	2259568	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.47	164	1391364	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.23	188	3035398	20.00	ng/ul	0.00
78) Chrysene-d12	21.33	240	3140859	20.00	ng/ul	0.00
86) Perylene-d12	23.69	264	3467139	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	27349	2.10	ng/uL	0.00
5) Phenol-d5	6.99	99	98173	2.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	183531	6.97	ng/ul	0.00
9) 2-Chlorophenol-d4	7.36	132	245778	6.41	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	170662	4.57	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	129852	6.64	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	148903	6.98	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	297944	7.53	ng/ul	0.00
29) 4-Chloroaniline-d4	10.75	131	302019	6.74	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	867414	7.79	ng/ul	0.00
47) Acenaphthylene-d8	14.16	160	1006374	7.28	ng/ul	0.00
52) 4-Nitrophenol-d4	14.67	143	28926	1.44	ng/ul	0.00
58) Fluorene-d10	15.47	176	679625	6.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.59	200	70745	3.61	ng/ul	0.00
71) Anthracene-d10	17.33	188	1101318	7.34	ng/ul	0.00
79) Pyrene-d10	19.57	212	1278860	7.62	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.53	264	1374607	7.20	ng/ul	0.00

Target Compounds

				Ovalue	
2) 1,4-Dioxane	3.35	88	512517	38.594	ng/uL# 96
6) Phenol	7.02	94	55856	1.183	ng/ul 97
45) Dimethylphthalate	13.92	163	166857	1.495	ng/ul 99

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

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