

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022220\
 Data File : BN009852.D
 Acq On : 22 Feb 2020 18:57
 Operator : CG/JU
 Sample : L1516-14
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBD72

Quant Time: Feb 24 02:42:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 19 04:46:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	109355	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.13	136	458865	20.00	ng/ul	-0.03
36) Acenaphthene-d10	14.03	164	285531	20.00	ng/ul	-0.02
62) Phenanthrene-d10	16.79	188	595735	20.00	ng/ul	-0.02
78) Chrysene-d12	21.01	240	532801	20.00	ng/ul	-0.01
86) Perylene-d12	23.10	264	593677	20.00	ng/ul	-0.03

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.03	96	17617	6.63	ng/uL	-0.02
5) Phenol-d5	6.59	99	332940	35.79	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	6.75	67	244553	38.10	ng/ul	-0.02
9) 2-Chlorophenol-d4	6.93	132	260592	37.21	ng/ul	-0.03
13) 4-Methylphenol-d8	8.10	113	264324	35.72	ng/ul	-0.02
19) Nitrobenzene-d5	8.53	128	127153	37.88	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.25	143	142154	38.71	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	9.78	165	248481	35.19	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.29	131	260874	32.89	ng/ul	-0.02
44) Dimethylphthalate-d6	13.46	166	794188	37.24	ng/ul	-0.02
47) Acenaphthylene-d8	13.72	160	957523	38.12	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.27	143	112330	28.84	ng/ul	-0.02
58) Fluorene-d10	15.03	176	694799	37.73	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.19	200	108519	29.32	ng/ul	-0.02
71) Anthracene-d10	16.89	188	1051891	38.52	ng/ul	-0.02
79) Pyrene-d10	19.19	212	1113852	40.00	ng/ul	-0.02
90) Benzo(a)pyrene-d12	22.97	264	1160725	39.15	ng/ul	-0.03

Target Compounds					Ovalue	
6) Phenol	6.62	94	12586	1.263	ng/ul	99
45) Dimethylphthalate	13.50	163	69284	3.113	ng/ul	97

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

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