

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP030220\  
 Data File : BP002006.D  
 Acq On : 03 Mar 2020 01:29  
 Operator : CG/JU  
 Sample : L1616-16  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBDR0

Quant Time: Mar 03 03:56:59 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP022720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 03 03:51:53 2020  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	297236	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1209817	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	763948	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1669129	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1652365	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1862911	20.00	ng/ul	0.00

## System Monitoring Compounds

3) 1,4-Dioxane-d8	3.17	96	30565	4.42	ng/uL	0.00
5) Phenol-d5	6.82	99	522484	23.04	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	311252	25.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	472081	24.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	400260	21.48	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	227100	24.44	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	231263	22.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	457565	22.72	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	469540	21.27	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1387046	23.99	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1734441	25.08	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	169607	15.35	ng/ul	0.00
58) Fluorene-d10	15.27	176	1276869	25.52	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	125154	13.60	ng/ul	0.00
71) Anthracene-d10	17.13	188	1922366	25.63	ng/ul	0.00
79) Pyrene-d10	19.37	212	2251842	26.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2392191	25.10	ng/ul	0.00

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.73	163	113777	1.906	ng/ul	99

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

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