

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002020.D
 Acq On : 03 Mar 2020 12:27
 Operator : CG/JU
 Sample : L1617-05
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDR9

Quant Time: Mar 04 06:01:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 03 05:31:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	265102	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1102909	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	706897	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1549805	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1513895	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1697716	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.16	96	35197	5.71	ng/uL	0.00
5) Phenol-d5	6.82	99	576150	28.49	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	334377	31.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	519493	29.70	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	469965	28.28	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	235318	27.78	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	259786	27.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	498206	27.13	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	181093	9.00	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1490054	27.85	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1840657	28.77	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	181857	17.78	ng/ul	0.00
58) Fluorene-d10	15.27	176	1363531	29.46	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	138096	16.16	ng/ul	0.00
71) Anthracene-d10	17.13	188	2026525	29.09	ng/ul	0.00
79) Pyrene-d10	19.37	212	2377570	30.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2558227	29.45	ng/ul	0.00
Target Compounds						
6) Phenol	6.85	94	29264	1.394	ng/ul	97
45) Dimethylphthalate	13.73	163	180142	3.262	ng/ul	100

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

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