

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030420\
 Data File : BP002062.D
 Acq On : 04 Mar 2020 21:58
 Operator : CG/JU
 Sample : L1677-12
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CC2

Quant Time: Mar 05 07:00:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 05 06:21:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	253781	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1040115	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	649184	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1386964	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1361126	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1533840	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.16	96	38558	6.53	ng/uL	0.00
5) Phenol-d5	6.82	99	689832	35.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.97	67	394558	38.49	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	610085	36.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	581724	36.57	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	288911	36.16	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	318528	35.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	624125	36.04	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	842871	44.40	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1887178	38.41	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2303888	39.21	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	241238	25.68	ng/ul	0.00
58) Fluorene-d10	15.27	176	1683389	39.60	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	178002	23.28	ng/ul	0.00
71) Anthracene-d10	17.13	188	2489348	39.93	ng/ul	0.00
79) Pyrene-d10	19.37	212	2871032	40.29	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3178125	40.50	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.73	163	521474	10.283	ng/ul	100

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

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