

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030620\
 Data File : BP002110.D
 Acq On : 06 Mar 2020 22:43
 Operator : CG/JU
 Sample : L1780-18
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0DJ7

Quant Time: Mar 07 02:20:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Mar 07 00:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	247531	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1022785	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	655404	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1431586	20.00	ng/ul	-0.01
78) Chrysene-d12	21.12	240	1430985	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1636047	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	32961	5.72	ng/uL	0.00
5) Phenol-d5	6.82	99	122911	6.51	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	331868	33.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	416822	25.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	251342	16.20	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	249142	31.71	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	264672	29.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	497461	29.21	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	37125	1.99	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1781127	35.91	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2098517	35.38	ng/ul	0.00
52) 4-Nitrophenol-d4	14.48	143	29097	3.07	ng/ul	0.00
58) Fluorene-d10	15.27	176	1578795	36.79	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	185670	23.53	ng/ul	0.00
71) Anthracene-d10	17.12	188	2523504	39.22	ng/ul	0.00
79) Pyrene-d10	19.37	212	2947442	39.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3325844	39.73	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.73	163	118700	2.318	ng/ul	99

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

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