

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
 Data File : BP001859.D
 Acq On : 24 Feb 2020 14:22
 Operator : CG/JU
 Sample : L1536-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB6R9

Quant Time: Feb 24 17:02:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 24 01:10:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	386436	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	1635039	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	1046239	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	2324030	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	2297052	20.00	ng/ul	0.00
86) Perylene-d12	23.35	264	2370706	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.18	96	61629	6.62	ng/uL	0.00
5) Phenol-d5	6.83	99	1072475	36.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	604669	35.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	960625	40.71	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	912800	39.41	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	478346	43.22	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	541161	54.06	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	961338	40.39	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	1308625	46.48	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	2986889	40.68	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	3588559	39.69	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	515126	40.29	ng/ul	0.00
58) Fluorene-d10	15.29	176	2592426	39.26	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	411794	44.94	ng/ul	0.00
71) Anthracene-d10	17.14	188	3921681	38.30	ng/ul	0.00
79) Pyrene-d10	19.39	212	4466641	37.89	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.21	264	4680854	40.88	ng/ul	0.00

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

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