

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
 Data File : BP002118.D
 Acq On : 09 Mar 2020 13:46
 Operator : CG/JU
 Sample : L1707-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB719

Quant Time: Mar 09 15:40:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 09 15:02:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	297588	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1231368	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	788265	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1703929	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1683260	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1925847	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	44937	6.49	ng/uL	0.00
5) Phenol-d5	6.82	99	740542	32.62	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	414194	34.46	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	651058	33.15	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	605362	32.45	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	301016	31.82	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	324938	30.53	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	638726	31.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.54	131	834002	37.11	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1920683	32.20	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2345836	32.88	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	232015	20.34	ng/ul	0.00
58) Fluorene-d10	15.27	176	1719983	33.32	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	194368	20.69	ng/ul	0.00
71) Anthracene-d10	17.12	188	2589181	33.81	ng/ul	0.00
79) Pyrene-d10	19.37	212	2963434	33.63	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3276666	33.25	ng/ul	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.73	163	168387	2.734	ng/ul	99

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

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