

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021721\
 Data File : BP004775.D
 Acq On : 17 Feb 2021 16:19
 Operator : CG/JU
 Sample : PB134658BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK58

Quant Time: Feb 17 17:50:36 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 15 12:16:19 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	42096	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	151790	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.41	164	84338	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.16	188	162910	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	162814	20.00	ng/ul	-0.01
86) Perylene-d12	23.55	264	176685	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.25	96	9507	8.73	ng/uL	0.00
5) Phenol-d5	6.93	99	145839	43.17	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	81901	46.67	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	125978	49.20	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	115252	43.36	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	56571	53.98	ng/ul	0.00
22) 2-Nitrophenol-d4	9.64	143	63174	54.93	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.18	165	110754	47.84	ng/ul	0.00
29) 4-Chloroaniline-d4	10.69	131	139954	54.16	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	296484	48.43	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	379702	50.98	ng/ul	0.00
52) 4-Nitrophenol-d4	14.61	143	37975	35.49	ng/ul	0.00
58) Fluorene-d10	15.40	176	248883	47.44	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	43780	44.46	ng/ul	0.00
71) Anthracene-d10	17.26	188	376201	52.36	ng/ul	0.00
79) Pyrene-d10	19.50	212	418730	54.94	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.40	264	475544	53.20	ng/ul	-0.01

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

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