

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021721\
 Data File : BP004805.D
 Acq On : 18 Feb 2021 11:48
 Operator : CG/JU
 Sample : PB134658BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK58

Quant Time: Feb 18 12:24:50 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	34612	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	128860	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.41	164	70863	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	143257	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	151540	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	164873	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.25	96	6126	6.84	ng/uL	0.00
5) Phenol-d5	6.95	99	105078	37.83	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.10	67	55882	38.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	89032	42.29	ng/ul	0.00
13) 4-Methylphenol-d8	8.48	113	81686	37.38	ng/ul	0.01
19) Nitrobenzene-d5	8.92	128	38817	43.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	44586	45.66	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	77100	39.23	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	98681	44.98	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	221533	43.07	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	275416	44.01	ng/ul	0.00
52) 4-Nitrophenol-d4	14.64	143	28860	32.10	ng/ul	0.03
58) Fluorene-d10	15.41	176	182819	41.48	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	27766	32.06	ng/ul	0.00
71) Anthracene-d10	17.27	188	284094	44.97	ng/ul	0.00
79) Pyrene-d10	19.51	212	318985	44.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.41	264	382754	45.89	ng/ul	0.00

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

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