

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020911.D
 Acq On : 11 Jul 2024 21:03
 Operator : MA/JU
 Sample : PB161883BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK883

Quant Time: Jul 11 23:51:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 10 18:38:19 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	211275	20.000	ng/u1	0.00
20) Naphthalene-d8	10.487	136	841460	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	535667	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.169	188	1166603	20.000	ng/u1	-0.01
79) Chrysene-d12	21.615	240	1172844	20.000	ng/u1	0.00
88) Perylene-d12	24.974	264	1383984	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	29365	6.325	ng/uL	0.00
4) Pyridine-d5	3.669	84	391101	29.097	ng/u1	0.00
7) Phenol-d5	6.922	99	491471	28.533	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	306962	29.407	ng/u1	0.00
11) 2-Chlorophenol-d4	7.269	132	418826	29.512	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	398813	27.877	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	199281	30.491	ng/u1	0.00
24) 2-Nitrophenol-d4	9.592	143	228152	30.499	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	383356	28.341	ng/u1	0.00
31) 4-Chloroaniline-d4	10.639	131	524560	29.273	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	1161040	29.816	ng/u1	-0.01
49) Acenaphthylene-d8	14.045	160	1308422	29.924	ng/u1	0.00
54) 4-Nitrophenol-d4	14.622	143	164509	29.693	ng/u1	0.01
60) Fluorene-d10	15.375	176	1005639	31.479	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.522	200	177020	25.079	ng/u1	0.01
73) Anthracene-d10	17.269	188	1627550	31.411	ng/u1	0.00
81) Pyrene-d10	19.651	212	1997934	27.554	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.745	264	2150918	31.512	ng/u1	-0.01

Target Compounds	Qvalue

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

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