

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
 Data File : BP018628.D
 Acq On : 06 Dec 2023 08:59
 Operator : MA/JU
 Sample : PB157406BL
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK406

Quant Time: Dec 06 21:38:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	259132	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.640	136	1094502	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.481	164	788519	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.228	188	1855715	20.000	ng/u1	0.00
79) Chrysene-d12	21.316	240	1622807	20.000	ng/u1	0.00
88) Perylene-d12	23.692	264	1622994	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	52567	6.913	ng/uL	-0.01
4) Pyridine-d5	3.675	84	788609	38.717	ng/u1	-0.01
7) Phenol-d5	7.005	99	641700	24.676	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.175	67	381579	24.757	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.369	132	488574	24.078	ng/u1	-0.02
15) 4-Methylphenol-d8	8.552	113	522509	24.828	ng/u1	-0.02
21) Nitrobenzene-d5	8.999	128	248183	25.477	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.722	143	276244	27.352	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.257	165	538057	26.054	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.781	131	776412	28.768	ng/u1	-0.01
46) Dimethylphthalate-d6	13.892	166	1913616	27.191	ng/u1	-0.01
49) Acenaphthylene-d8	14.169	160	1998441	25.788	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.675	143	297405	25.745	ng/u1	-0.01
60) Fluorene-d10	15.475	176	1584362	26.820	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	248374	22.322	ng/u1	0.00
73) Anthracene-d10	17.328	188	2625471	26.912	ng/u1	0.00
81) Pyrene-d10	19.569	212	3096408	24.512	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	2578131	27.244	ng/u1	0.00

Target Compounds	Qvalue

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

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