

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017883.D
 Acq On : 02 Nov 2023 13:17
 Operator : MA/JU
 Sample : PB156594BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK594

Quant Time: Nov 02 23:36:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	145632	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.704	136	548483	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.575	164	316557	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.369	188	635488	20.000	ng/u1	0.00
79) Chrysene-d12	21.516	240	529198	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	542447	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	22854	5.838	ng/uL	0.01
4) Pyridine-d5	3.699	84	270101	26.353	ng/u1	0.00
7) Phenol-d5	7.040	99	396177	32.691	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	245906	32.856	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.399	132	332224	33.611	ng/u1	-0.01
15) 4-Methylphenol-d8	8.599	113	317086	33.319	ng/u1	-0.01
21) Nitrobenzene-d5	9.081	128	157846	35.516	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.799	143	178964	36.444	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	308517	32.028	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.881	131	397295	31.515	ng/u1	-0.01
46) Dimethylphthalate-d6	13.992	166	894908	33.676	ng/u1	-0.01
49) Acenaphthylene-d8	14.269	160	1012584	34.168	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.828	143	106087	28.017	ng/u1	0.00
60) Fluorene-d10	15.581	176	737070	32.087	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	117367	28.039	ng/u1	0.00
73) Anthracene-d10	17.469	188	1144716	35.249	ng/u1	0.00
81) Pyrene-d10	19.733	212	1280391	38.591	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	1069123	35.387	ng/u1	0.00

Target Compounds Qvalue

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

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