

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018517.D
 Acq On : 02 Dec 2023 16:02
 Operator : MA/JU
 Sample : PB157321BL
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK321

Quant Time: Dec 02 21:32:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 01 21:56:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	401208	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	1696149	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.486	164	1041333	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	2135516	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	2024084	20.000	ng/ul	-0.01
88) Perylene-d12	23.692	264	2412607	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	68860	5.849	ng/uL	0.00
4) Pyridine-d5	3.687	84	853271	27.057	ng/ul	0.00
7) Phenol-d5	7.028	99	1155222	28.692	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	706805	29.619	ng/ul	0.00
11) 2-Chlorophenol-d4	7.393	132	891263	28.369	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	911617	27.978	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128	435590	28.854	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	462921	29.577	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	873897	27.306	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.798	131	994740	23.783	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	2645435	28.464	ng/ul	0.00
49) Acenaphthylene-d8	14.181	160	3024891	29.557	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.686	143	433864	28.439	ng/ul	0.00
60) Fluorene-d10	15.481	176	2231598	28.605	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	307273	23.997	ng/ul	0.00
73) Anthracene-d10	17.333	188	3383841	30.141	ng/ul	0.00
81) Pyrene-d10	19.569	212	3897713	24.738	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	4211991	29.942	ng/ul	-0.01

Target Compounds Qvalue

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

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