

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP023003.D
 Acq On : 10 Nov 2024 16:55
 Operator : RC/JU
 Sample : PB164817BL
 Misc :
 ALS Vial : 94 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK817

Quant Time: Nov 10 22:41:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 08 23:49:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	94802	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.334	136	371019	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.204	164	308986	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.022	188	760031	20.000	ng/ul	-0.01
79) Chrysene-d12	21.457	240	853155	20.000	ng/ul	0.00
88) Perylene-d12	24.686	264	963772	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	11039	5.597	ng/uL	0.00
4) Pyridine-d5	3.552	84	153031	26.116	ng/ul	0.00
7) Phenol-d5	6.793	99	189657	27.059	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	120427	29.245	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	148294	28.889	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	158164	27.422	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	76895	29.770	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	93175	30.258	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	172562	26.759	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.487	131	206830	26.885	ng/ul	-0.01
46) Dimethylphthalate-d6	13.610	166	693461	31.139	ng/ul	0.00
49) Acenaphthylene-d8	13.887	160	693893	29.862	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.469	143	84574	25.271	ng/ul	0.00
60) Fluorene-d10	15.222	176	582882	29.853	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.428	200	1760	0.387	ng/ul	0.06
73) Anthracene-d10	17.110	188	1033576	32.118	ng/ul	-0.01
81) Pyrene-d10	19.516	212	1306122	33.237	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.463	264	1465855	31.930	ng/ul	0.00

Target Compounds	Qvalue

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

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