

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
 Data File : BP022171.D
 Acq On : 02 Oct 2024 21:34
 Operator : RC/JU
 Sample : PB163489BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK489

Quant Time: Oct 03 07:34:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	120575	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.457	136	465120	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.310	164	374188	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.116	188	875804	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.539	240	922193	20.000	ng/u1	0.00
88) Perylene-d12	24.809	264	1028353	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	12912	4.195	ng/uL	0.00
4) Pyridine-d5	3.634	84	188639	21.439	ng/u1	0.00
7) Phenol-d5	6.881	99	215187	21.817	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.034	67	132203	21.768	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.240	132	176475	24.730	ng/u1	0.00
15) 4-Methylphenol-d8	8.399	113	179973	23.027	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	85017	24.421	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.546	143	97415	24.679	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.087	165	194100	22.643	ng/u1	0.00
31) 4-Chloroaniline-d4	10.593	131	237900	23.260	ng/u1	-0.01
46) Dimethylphthalate-d6	13.716	166	692442	23.939	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	726466	23.722	ng/u1	0.00
54) 4-Nitrophenol-d4	14.522	143	89602	18.110	ng/u1	0.00
60) Fluorene-d10	15.322	176	585743	22.893	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	98540	18.123	ng/u1	0.00
73) Anthracene-d10	17.204	188	980462	23.980	ng/u1	-0.01
81) Pyrene-d10	19.592	212	1184286	24.788	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.586	264	1256097	23.223	ng/u1	-0.02

Target Compounds Qvalue

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

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