

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023636.D
 Acq On : 17 Jan 2025 20:42
 Operator : RC/JU
 Sample : PB166106BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK106

Quant Time: Jan 19 23:07:28 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 14 16:27:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	496639	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1994374	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1272020	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2628726	20.000	ng/u1	0.00
79) Chrysene-d12	21.692	240	2594858	20.000	ng/u1	-0.02
88) Perylene-d12	25.121	264	2857389	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	89360	6.634	ng/uL	0.00
4) Pyridine-d5	3.716	84	1295738	34.507	ng/u1	0.00
7) Phenol-d5	6.963	99	1508241	35.275	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	832415	33.233	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	1201960	35.596	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	1167698	35.209	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	527734	34.668	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	408832	40.917	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	1051531	34.599	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	1620258	33.394	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	3365320	33.995	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	3792046	32.200	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	476462	28.955	ng/u1	0.00
60) Fluorene-d10	15.445	176	2885136	33.675	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	203058	23.741	ng/u1	0.00
73) Anthracene-d10	17.333	188	4618769	32.549	ng/u1	-0.01
81) Pyrene-d10	19.698	212	5289496	36.796	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.892	264	5247899	33.796	ng/u1	-0.02

Target Compounds Qvalue

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

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