

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012725\
 Data File : BP023773.D
 Acq On : 28 Jan 2025 15:00
 Operator : RC/JU
 Sample : PB166284BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK284

Quant Time: Jan 29 08:31:39 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012225.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 27 21:53:59 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	434088	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.558	136	1755756	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.410	164	1159817	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.210	188	2470440	20.000	ng/u1	0.00
79) Chrysene-d12	21.669	240	2258805	20.000	ng/u1	0.00
88) Perylene-d12	25.080	264	2481677	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	64729	6.006	ng/uL	-0.01
4) Pyridine-d5	3.699	84	867940	27.907	ng/u1	0.00
7) Phenol-d5	6.964	99	1179664	32.116	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.111	67	578906	27.903	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.317	132	970241	34.746	ng/u1	-0.01
15) 4-Methylphenol-d8	8.487	113	935890	32.403	ng/u1	0.01
21) Nitrobenzene-d5	8.922	128	427427	32.311	ng/u1	0.00
24) 2-Nitrophenol-d4	9.640	143	449152	37.035	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	902865	35.418	ng/u1	0.01
31) 4-Chloroaniline-d4	10.687	131	1106401	27.139	ng/u1	0.00
46) Dimethylphthalate-d6	13.816	166	2707730	32.462	ng/u1	0.00
49) Acenaphthylene-d8	14.099	160	3064175	31.741	ng/u1	0.00
54) 4-Nitrophenol-d4	14.640	143	447150	26.680	ng/u1	0.05
60) Fluorene-d10	15.422	176	2400467	33.620	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.534	200	294931	24.243	ng/u1	0.00
73) Anthracene-d10	17.316	188	3826356	32.262	ng/u1	0.00
81) Pyrene-d10	19.686	212	4023977	35.010	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.845	264	4277025	34.432	ng/u1	0.00

Target Compounds Qvalue

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

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