

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012725\
 Data File : BP023749.D
 Acq On : 27 Jan 2025 21:16
 Operator : RC/JU
 Sample : PB166248BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK248

Quant Time: Jan 27 23:28:40 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	7.787	152	457543	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	1851043	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	1242807	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.228	188	2657302	20.000	ng/u1	0.01
79) Chrysene-d12	21.669	240	2845516	20.000	ng/u1	0.00
88) Perylene-d12	25.074	264	3069158	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	77503	6.823	ng/uL	0.00
4) Pyridine-d5	3.711	84	1086083	33.131	ng/u1	0.00
7) Phenol-d5	6.975	99	1416842	36.596	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.122	67	708769	32.411	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	1142127	38.805	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	1143498	37.561	ng/u1	0.02
21) Nitrobenzene-d5	8.934	128	512816	36.771	ng/u1	0.00
24) 2-Nitrophenol-d4	9.652	143	531939	41.604	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	1067757	39.730	ng/u1	0.02
31) 4-Chloroaniline-d4	10.705	131	1216855	28.312	ng/u1	0.01
46) Dimethylphthalate-d6	13.822	166	3292582	36.838	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	3682508	35.599	ng/u1	0.01
54) 4-Nitrophenol-d4	14.640	143	592572	32.996	ng/u1	0.05
60) Fluorene-d10	15.434	176	2823762	36.907	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.545	200	387185	29.588	ng/u1	0.02
73) Anthracene-d10	17.328	188	4649784	36.448	ng/u1	0.00
81) Pyrene-d10	19.698	212	5487067	37.896	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.833	264	5867090	38.192	ng/u1	-0.01

Target Compounds	Qvalue

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

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