

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
 Data File : BP023985.D
 Acq On : 07 Feb 2025 01:28
 Operator : RC/JU
 Sample : PB166547BL
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK547

Quant Time: Feb 07 08:00:23 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 11:19:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	450061	20.000	ng/u1	0.00
20) Naphthalene-d8	10.540	136	1924756	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.387	164	1240655	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	2453139	20.000	ng/u1	0.00
79) Chrysene-d12	21.627	240	2230567	20.000	ng/u1	-0.01
88) Perylene-d12	25.004	264	2236414	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	87487	8.077	ng/uL	0.00
4) Pyridine-d5	3.699	84	1245998	39.864	ng/u1	0.00
7) Phenol-d5	6.952	99	1740488	43.699	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	869459	41.242	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	1408035	45.019	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	1381834	42.471	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	640484	41.592	ng/u1	0.00
24) 2-Nitrophenol-d4	9.628	143	686835	41.335	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	1295089	42.486	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1832449	40.197	ng/u1	0.00
46) Dimethylphthalate-d6	13.793	166	3765234	40.687	ng/u1	0.00
49) Acenaphthylene-d8	14.081	160	4369236	41.862	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	595631	32.410	ng/u1	0.00
60) Fluorene-d10	15.392	176	3309065	42.462	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	326204	21.554	ng/u1	0.00
73) Anthracene-d10	17.281	188	5205374	43.216	ng/u1	0.00
81) Pyrene-d10	19.651	212	5760323	48.196	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.768	264	5212256	44.669	ng/u1	0.00

Target Compounds Qvalue

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

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