

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
 Data File : BP023986.D
 Acq On : 07 Feb 2025 02:08
 Operator : RC/JU
 Sample : PB166550BL
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK550

Quant Time: Feb 07 07:34:48 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	491359	20.000	ng/u1	0.00
20) Naphthalene-d8	10.546	136	2100497	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.398	164	1327023	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.204	188	2621940	20.000	ng/u1	0.02
79) Chrysene-d12	21.639	240	2469483	20.000	ng/u1	0.00
88) Perylene-d12	25.016	264	2643970	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	84281	7.127	ng/uL	0.00
4) Pyridine-d5	3.699	84	1239310	36.317	ng/u1	0.00
7) Phenol-d5	6.952	99	1639033	37.693	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	818404	35.557	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	1320055	38.658	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	1288823	36.283	ng/u1	0.00
21) Nitrobenzene-d5	8.911	128	592771	35.273	ng/u1	0.00
24) 2-Nitrophenol-d4	9.628	143	645615	35.604	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	1196787	35.977	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1713724	34.447	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	3379962	34.147	ng/u1	0.01
49) Acenaphthylene-d8	14.087	160	4012206	35.939	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	548959	27.926	ng/u1	0.00
60) Fluorene-d10	15.410	176	2978883	35.737	ng/u1	0.02
65) 4,6-Dinitro-2-methylph...	15.522	200	251573	15.553	ng/u1	0.01
73) Anthracene-d10	17.292	188	4840567	37.600	ng/u1	0.00
81) Pyrene-d10	19.669	212	5342212	40.374	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.780	264	5239887	37.984	ng/u1	0.00

Target Compounds Qvalue

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

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