

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
 Data File : BP023998.D
 Acq On : 07 Feb 2025 11:47
 Operator : RC/JU
 Sample : PB166554BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK554

Quant Time: Feb 12 11:34:05 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	385093	20.000	ng/u1	0.00
20) Naphthalene-d8	10.540	136	1714903	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.387	164	1098313	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.175	188	2150024	20.000	ng/u1	-0.01
79) Chrysene-d12	21.627	240	1929205	20.000	ng/u1	-0.01
88) Perylene-d12	24.986	264	1970599	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	58219	6.281	ng/uL	0.00
4) Pyridine-d5	3.699	84	881568	32.962	ng/u1	0.00
7) Phenol-d5	6.952	99	1173895	34.446	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	585904	32.480	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	940979	35.161	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	924105	33.194	ng/u1	0.00
21) Nitrobenzene-d5	8.911	128	424492	30.939	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	455668	30.779	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	857927	31.589	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	1228418	30.244	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2474689	30.207	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	2948660	31.913	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	352348	21.657	ng/u1	0.00
60) Fluorene-d10	15.393	176	2204065	31.948	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	121114	9.131	ng/u1	0.00
73) Anthracene-d10	17.281	188	3459419	32.770	ng/u1	0.00
81) Pyrene-d10	19.657	212	3747070	36.249	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.763	264	3409236	33.158	ng/u1	-0.01

Target Compounds Qvalue

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

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