

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012025\
 Data File : BP023676.D
 Acq On : 21 Jan 2025 10:05
 Operator : RC/JU
 Sample : PB166118BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK118

Quant Time: Jan 21 22:19:33 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 20 21:50:54 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	852443	20.000	ng/u1	-0.04
20) Naphthalene-d8	10.552	136	3386921	20.000	ng/u1	-0.03
38) Acenaphthene-d10	14.434	164	2143956	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.245	188	4231909	20.000	ng/u1	0.02
79) Chrysene-d12	21.686	240	4326586	20.000	ng/u1	0.02
88) Perylene-d12	25.080	264	4663215	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	136767	5.916	ng/uL	-0.02
4) Pyridine-d5	3.693	84	2098696	32.562	ng/u1	-0.02
7) Phenol-d5	6.922	99	2458754	33.503	ng/u1	-0.04
9) Bis-(2-Chloroethyl)eth...	7.093	67	1331813	30.978	ng/u1	-0.05
11) 2-Chlorophenol-d4	7.299	132	1958977	33.800	ng/u1	-0.04
15) 4-Methylphenol-d8	8.452	113	1888836	33.181	ng/u1	-0.04
21) Nitrobenzene-d5	8.910	128	879941	34.038	ng/u1	-0.04
24) 2-Nitrophenol-d4	9.634	143	841379	49.585	ng/u1	-0.03
28) 2,4-Dichlorophenol-d3	10.163	165	1683155	32.611	ng/u1	-0.04
31) 4-Chloroaniline-d4	10.681	131	2344608	28.455	ng/u1	-0.02
46) Dimethylphthalate-d6	13.828	166	5051276	30.274	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	5822238	29.332	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	882103	31.805	ng/u1	0.01
60) Fluorene-d10	15.434	176	4348097	30.111	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	460744	33.461	ng/u1	0.00
73) Anthracene-d10	17.345	188	6839206	29.938	ng/u1	0.02
81) Pyrene-d10	19.710	212	7960285	33.211	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.857	264	8139262	32.118	ng/u1	0.05

Target Compounds Qvalue

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

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