

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012025\
 Data File : BP023653.D
 Acq On : 20 Jan 2025 13:52
 Operator : RC/JU
 Sample : PB166120BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK120

Quant Time: Jan 20 22:54:38 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 14 16:27:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	699314	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	2769782	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1660891	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.222	188	3302231	20.000	ng/u1	-0.02
79) Chrysene-d12	21.674	240	3289627	20.000	ng/u1	-0.04
88) Perylene-d12	25.068	264	3659722	20.000	ng/u1	-0.08
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	118489	6.247	ng/uL	0.00
4) Pyridine-d5	3.716	84	1789381	33.842	ng/u1	0.00
7) Phenol-d5	6.963	99	2062838	34.263	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	1153120	32.695	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	1674921	35.227	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	1593996	34.133	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	752865	35.612	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	680556	49.044	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	1395704	33.067	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	2141165	31.776	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	4178075	32.323	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	4944509	32.155	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	666847	31.037	ng/u1	0.00
60) Fluorene-d10	15.439	176	3626988	32.422	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	341905	31.821	ng/u1	0.00
73) Anthracene-d10	17.322	188	5626260	31.562	ng/u1	-0.02
81) Pyrene-d10	19.692	212	6398312	35.109	ng/u1	-0.03
92) Benzo(a)pyrene-d12	24.827	264	6797914	34.180	ng/u1	-0.08

Target Compounds Qvalue

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

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