

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
 Data File : BP023931.D
 Acq On : 05 Feb 2025 10:50
 Operator : RC/JU
 Sample : PB166502BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK502

Quant Time: Feb 05 11:12:42 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	7.775	152	318095	20.000	ng/u1	0.00
20) Naphthalene-d8	10.546	136	1317461	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.387	164	806656	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	1671879	20.000	ng/u1	0.00
79) Chrysene-d12	21.622	240	1689612	20.000	ng/u1	-0.02
88) Perylene-d12	24.980	264	1733053	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	59865	7.819	ng/uL	0.00
4) Pyridine-d5	3.705	84	854088	38.661	ng/u1	0.00
7) Phenol-d5	6.958	99	1027634	36.505	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	548470	36.809	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	826923	37.407	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	800387	34.806	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	370989	35.196	ng/u1	0.00
24) 2-Nitrophenol-d4	9.634	143	378225	33.255	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	725822	34.787	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1072852	34.383	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2127805	35.364	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	2536415	37.376	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	358011	29.961	ng/u1	-0.01
60) Fluorene-d10	15.392	176	1938001	38.248	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	211551	20.511	ng/u1	0.00
73) Anthracene-d10	17.286	188	3144300	38.303	ng/u1	0.00
81) Pyrene-d10	19.663	212	3688553	40.743	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.757	264	3580239	39.594	ng/u1	-0.02

Target Compounds Qvalue

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

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