

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023498.D
 Acq On : 18 Dec 2024 15:00
 Operator : RC/JU
 Sample : PB165651BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK651

Quant Time: Dec 18 15:24:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.505	152	292475	20.000	ng/u1	-0.04
20) Naphthalene-d8	10.251	136	1097336	20.000	ng/u1	#-0.03
38) Acenaphthene-d10	14.134	164	709439	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.951	188	1625155	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	1780774	20.000	ng/u1	# 0.00
88) Perylene-d12	24.568	264	1794140	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	37119	4.672	ng/uL	-0.02
4) Pyridine-d5	3.493	84	536547	26.265	ng/u1	-0.02
7) Phenol-d5	6.722	99	610578	26.607	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	6.852	67	381401	26.343	ng/u1	-0.04
11) 2-Chlorophenol-d4	7.057	132	437836	25.851	ng/u1	-0.03
15) 4-Methylphenol-d8	8.222	113	463772	25.040	ng/u1	-0.04
21) Nitrobenzene-d5	8.652	128	197385	24.088	ng/u1	-0.04
24) 2-Nitrophenol-d4	9.363	143	236043	24.994	ng/u1	-0.03
28) 2,4-Dichlorophenol-d3	9.910	165	448462	22.500	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.410	131	571373	25.589	ng/u1	-0.03
46) Dimethylphthalate-d6	13.545	166	1449063	26.311	ng/u1	-0.01
49) Acenaphthylene-d8	13.822	160	1486322	25.586	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.439	143	195642	28.989	ng/u1	0.03
60) Fluorene-d10	15.157	176	1190374	24.573	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.316	200	225717	22.981	ng/u1	0.00
73) Anthracene-d10	17.057	188	1990939	26.843	ng/u1	0.00
81) Pyrene-d10	19.445	212	2404180	25.841	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.357	264	2411445	25.108	ng/u1	0.01

Target Compounds Qvalue

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

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