

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112524\
 Data File : BP023312.D
 Acq On : 25 Nov 2024 12:18
 Operator : RC/JU
 Sample : PB165168BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK168

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/27/2024
 Supervised By :mohammad ahmed 11/27/2024

Quant Time: Nov 26 00:00:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 04:34:35 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.546	152	68607	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.293	136	263744	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.163	164	218141	20.000	ng/u1	-0.02
64) Phenanthrene-d10	16.981	188	566401	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.410	240	657443	20.000	ng/u1	-0.01
88) Perylene-d12	24.610	264	722144	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.117	96	9535	6.680	ng/uL	-0.02
4) Pyridine-d5	3.534	84	122541	28.897	ng/u1	-0.01
7) Phenol-d5	6.758	99	140365	27.673	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	6.893	67	95451	32.030	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.099	132	107147	28.843	ng/u1	-0.02
15) 4-Methylphenol-d8	8.258	113	112759	27.014	ng/u1	-0.02
21) Nitrobenzene-d5	8.693	128	53876	29.342	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.405	143	66163	30.226	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	9.957	165	120670	26.324	ng/u1	0.00
31) 4-Chloroaniline-d4	10.457	131	147347m	26.943	ng/u1	-0.01
46) Dimethylphthalate-d6	13.569	166	514515	32.726	ng/u1	-0.02
49) Acenaphthylene-d8	13.857	160	507215	30.919	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.463	143	59805m	25.312	ng/u1	0.00
60) Fluorene-d10	15.192	176	442074	32.071	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.339	200	80152	23.623	ng/u1	-0.01
73) Anthracene-d10	17.075	188	752169	31.364	ng/u1	-0.02
81) Pyrene-d10	19.463	212	1014212	33.492	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.374	264	1150174	33.437	ng/u1	-0.05

Target Compounds Qvalue

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

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