

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
 Data File : BP020926.D
 Acq On : 12 Jul 2024 08:35
 Operator : MA/JU
 Sample : PB161865BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK865

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/12/2024
 Supervised By :mohammad ahmed 07/13/2024

Quant Time: Jul 12 09:31:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 10 18:38:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	294040	20.000	ng/u1	0.00
20) Naphthalene-d8	10.481	136	1219396	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.357	164	713098	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.169	188	1285528	20.000	ng/u1	-0.01
79) Chrysene-d12	21.621	240	1285835	20.000	ng/u1	0.00
88) Perylene-d12	24.974	264	1496714	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	35096	5.432	ng/uL	0.00
4) Pyridine-d5	3.669	84	475851	25.437	ng/u1	0.00
7) Phenol-d5	6.916	99	606349	25.294	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	382658	26.340	ng/u1	0.00
11) 2-Chlorophenol-d4	7.263	132	509174	25.780	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	499163	25.071	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	252463	26.656	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	290600	26.807	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	469779	23.966	ng/u1	0.00
31) 4-Chloroaniline-d4	10.640	131	661603	25.478	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	1377384	26.571	ng/u1	-0.01
49) Acenaphthylene-d8	14.039	160	1524746	26.195	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.622	143	159404m	21.613	ng/u1	0.01
60) Fluorene-d10	15.369	176	1060907	24.946	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	153043	19.676	ng/u1	0.00
73) Anthracene-d10	17.275	188	1552928	27.198	ng/u1	0.00
81) Pyrene-d10	19.657	212	1855243	23.338	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.739	264	2036757	27.592	ng/u1	-0.02

Target Compounds Qvalue

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

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