

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
 Data File : BP021141.D
 Acq On : 25 Jul 2024 10:18
 Operator : MA/JU
 Sample : PB162100BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK100

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/26/2024
 Supervised By :mohammad ahmed 07/27/2024

Quant Time: Jul 25 22:48:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 25 00:15:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	235597	20.000	ng/u1	0.00
20) Naphthalene-d8	10.405	136	904681	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.287	164	568987	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1201546	20.000	ng/u1	0.00
79) Chrysene-d12	21.563	240	1034594	20.000	ng/u1	0.00
88) Perylene-d12	24.886	264	1104497	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	27141	5.242	ng/uL	-0.01
4) Pyridine-d5	3.581	84	297495	19.848	ng/u1	-0.02
7) Phenol-d5	6.846	99	448231	23.336	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	6.987	67	275871	23.700	ng/u1	0.00
11) 2-Chlorophenol-d4	7.187	132	379676	23.992	ng/u1	-0.02
15) 4-Methylphenol-d8	8.363	113	362917	22.749	ng/u1	-0.02
21) Nitrobenzene-d5	8.799	128	179794	25.587	ng/u1	0.00
24) 2-Nitrophenol-d4	9.510	143	213971	26.605	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.057	165	344475	23.687	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.569	131	441039	22.892	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	1060477	25.639	ng/u1	0.00
49) Acenaphthylene-d8	13.975	160	1191698	25.659	ng/u1	0.00
54) 4-Nitrophenol-d4	14.593	143	142936m	24.288	ng/u1	0.00
60) Fluorene-d10	15.304	176	897731	26.456	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	138892	19.105	ng/u1	0.00
73) Anthracene-d10	17.210	188	1428241	26.763	ng/u1	0.00
81) Pyrene-d10	19.598	212	1689371	26.412	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.657	264	1467755	26.944	ng/u1	0.00

Target Compounds Qvalue

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

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