

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090324\
 Data File : BP021792.D
 Acq On : 03 Sep 2024 17:47
 Operator : RC/JU
 Sample : PB163070BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK070

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 09/04/2024
 Supervised By :mohammad ahmed 09/07/2024

Quant Time: Sep 04 04:42:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 29 23:56:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	315590	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.563	136	1246125	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	791337	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.240	188	1747453	20.000	ng/u1	0.01
79) Chrysene-d12	21.680	240	1792967	20.000	ng/u1	0.01
88) Perylene-d12	25.092	264	1959725	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	50834	5.348	ng/uL	0.00
4) Pyridine-d5	3.681	84	653871	23.781	ng/u1	0.00
7) Phenol-d5	6.952	99	823499	24.451	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	465564	25.401	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	542123	24.932	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	602692	23.158	ng/u1	-0.01
21) Nitrobenzene-d5	8.928	128	260911	26.025	ng/u1	0.00
24) 2-Nitrophenol-d4	9.646	143	266242	26.162	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	475046m	23.688	ng/u1	0.00
31) 4-Chloroaniline-d4	10.699	131	685701	23.502	ng/u1	0.00
46) Dimethylphthalate-d6	13.822	166	1643254	27.138	ng/u1	-0.02
49) Acenaphthylene-d8	14.110	160	1740276	25.599	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.628	143	254746m	22.114	ng/u1	0.00
60) Fluorene-d10	15.434	176	1372366	26.959	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.563	200	199590	20.767	ng/u1	0.00
73) Anthracene-d10	17.340	188	2209716	26.645	ng/u1	0.00
81) Pyrene-d10	19.704	212	2758140	26.910	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.851	264	2758714	26.399	ng/u1	0.01

Target Compounds Qvalue

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

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