

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090324\
 Data File : BP021806.D
 Acq On : 04 Sep 2024 04:29
 Operator : RC/JU
 Sample : PB163035BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK035

Manual Integrations
 APPROVED

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

Quant Time: Sep 04 06:05:02 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	7.781	152	361073	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	1405329	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.416	164	798091	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.210	188	1694130	20.000	ng/u1	-0.02
79) Chrysene-d12	21.663	240	1691119	20.000	ng/u1	0.00
88) Perylene-d12	25.068	264	1911548	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	52304	4.809	ng/uL	0.00
4) Pyridine-d5	3.687	84	695295	22.103	ng/u1	0.00
7) Phenol-d5	6.952	99	899491	23.343	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	508887	24.267	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	616911	24.798	ng/u1	0.00
15) 4-Methylphenol-d8	8.481	113	662030	22.233	ng/u1	0.00
21) Nitrobenzene-d5	8.928	128	292972	25.912	ng/u1	0.00
24) 2-Nitrophenol-d4	9.646	143	292324	25.471	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	545801m	24.133	ng/u1	0.00
31) 4-Chloroaniline-d4	10.693	131	764236	23.227	ng/u1	0.00
46) Dimethylphthalate-d6	13.822	166	1590365	26.043	ng/u1	-0.02
49) Acenaphthylene-d8	14.110	160	1751834	25.551	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.622	143	254433	21.899	ng/u1	0.00
60) Fluorene-d10	15.422	176	1347296	26.242	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.545	200	190765	20.473	ng/u1	-0.01
73) Anthracene-d10	17.316	188	2145247	26.682	ng/u1	-0.02
81) Pyrene-d10	19.680	212	2572646	26.612	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.833	264	2667398	26.169	ng/u1	0.00

Target Compounds Qvalue

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

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