

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090324\
 Data File : BP021807.D
 Acq On : 04 Sep 2024 05:12
 Operator : RC/JU
 Sample : P3733-09
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44A3

Manual Integrations
 APPROVED

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

Quant Time: Sep 04 06:08:37 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	222821	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	848680	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	490783	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.216	188	1040724	20.000	ng/u1	-0.01
79) Chrysene-d12	21.674	240	1181249	20.000	ng/u1	0.00
88) Perylene-d12	25.062	264	1377177	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	50547	7.531	ng/uL	0.00
4) Pyridine-d5	3.681	84	695021	35.802	ng/u1	0.00
7) Phenol-d5	6.952	99	863761	36.324	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	480233	37.109	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	593059	38.630	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	654411	35.614	ng/u1	-0.01
21) Nitrobenzene-d5	8.928	128	276604	40.510	ng/u1	0.00
24) 2-Nitrophenol-d4	9.646	143	277718	40.070	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	506301m	37.070	ng/u1	0.00
31) 4-Chloroaniline-d4	10.693	131	730208	36.749	ng/u1	0.00
46) Dimethylphthalate-d6	13.822	166	1603062	42.688	ng/u1	-0.02
49) Acenaphthylene-d8	14.110	160	1805401	42.820	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.622	143	269796	37.762	ng/u1	0.00
60) Fluorene-d10	15.428	176	1378748	43.670	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.545	200	191017	33.371	ng/u1	-0.01
73) Anthracene-d10	17.322	188	2203329	44.609	ng/u1	-0.02
81) Pyrene-d10	19.698	212	2855513	42.288	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.845	264	3396319	46.249	ng/u1	0.00

Target Compounds Qvalue

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

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