

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100324\
 Data File : BM047829.D
 Acq On : 04 Oct 2024 05:09
 Operator : RC/JU
 Sample : P4084-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCYM0

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/04/2024
 Supervised By :mohammad ahmed 10/06/2024

Quant Time: Oct 04 05:44:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	263260	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.592	136	1057650	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.433	164	656649	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.174	188	1314380	20.000	ng/u1	-0.01
79) Chrysene-d12	21.397	240	1218971	20.000	ng/u1	-0.01
88) Perylene-d12	24.397	264	1420729	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	19397	2.381	ng/uL	0.00
4) Pyridine-d5	3.669	84	172648	7.238	ng/u1	0.00
7) Phenol-d5	6.969	99	335195	13.465	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	215815	12.217	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.334	132	280136	15.467	ng/u1	-0.01
15) 4-Methylphenol-d8	8.504	113	253011	13.670	ng/u1	-0.01
21) Nitrobenzene-d5	8.957	128	136116m	16.164	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.675	143	142616	17.673	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.216	165	261787	16.239	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.728	131	385512	15.468	ng/u1	-0.01
46) Dimethylphthalate-d6	13.845	166	860463	18.074	ng/u1	-0.01
49) Acenaphthylene-d8	14.127	160	978090	17.328	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.627	143	110345	11.641	ng/u1	0.00
60) Fluorene-d10	15.427	176	747270	18.122	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.539	200	80630	10.483	ng/u1	-0.01
73) Anthracene-d10	17.268	188	1145932	17.731	ng/u1	-0.02
81) Pyrene-d10	19.556	212	1329369	19.215	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.186	264	1358299	18.427	ng/u1	-0.02

Target Compounds Qvalue

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

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