

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101124\
 Data File : BN034562.D
 Acq On : 11 Oct 2024 19:23
 Operator : RC/JU
 Sample : P4386-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DSN002

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/14/2024
 Supervised By :mohammad ahmed 10/15/2024

Quant Time: Oct 14 01:09:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 00:21:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.365	152	3574	0.400	ng	0.00
7) Naphthalene-d8	10.116	136	10243	0.400	ng	-0.01
13) Acenaphthene-d10	14.008	164	5664	0.400	ng	0.00
19) Phenanthrene-d10	16.771	188	10430	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	3567	0.400	ng	0.00
35) Perylene-d12	23.104	264	10493	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	2214	0.218	ng	0.00
5) Phenol-d6	6.578	99	1560m	0.132	ng	0.00
8) Nitrobenzene-d5	8.504	82	2339	0.299	ng	-0.01
11) 2-Methylnaphthalene-d10	11.716	152	5576	0.369	ng	-0.01
14) 2,4,6-Tribromophenol	15.517	330	985	0.384	ng	-0.02
15) 2-Fluorobiphenyl	12.629	172	7573	0.329	ng	0.00
27) Fluoranthene-d10	18.820	212	11485	0.436	ng	0.00
31) Terphenyl-d14	19.438	244	7826	1.099	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.018	88	2720	0.609	ng	# 83

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

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