

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034546.D
 Acq On : 11 Oct 2024 07:27
 Operator : RC/JU
 Sample : P4231-01DL 2X
 Misc :
 ALS Vial : 122 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D-20240924DL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/11/2024
 Supervised By :mohammad ahmed 10/14/2024

Quant Time: Oct 11 07:57:28 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	7.373	152	3893	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	11437	0.400	ng	0.01
13) Acenaphthene-d10	14.030	164	6850	0.400	ng	0.01
19) Phenanthrene-d10	16.783	188	12995	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	7759	0.400	ng	0.00
35) Perylene-d12	23.116	264	9721	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	558	0.051	ng	0.00
5) Phenol-d6	6.643	99	242m	0.019	ng	0.07
8) Nitrobenzene-d5	8.557	82	958m	0.110	ng	0.04
11) 2-Methylnaphthalene-d10	11.750	152	2412m	0.143	ng	0.02
14) 2,4,6-Tribromophenol	15.555	330	310	0.100	ng	0.02
15) 2-Fluorobiphenyl	12.651	172	4063	0.146	ng	0.01
27) Fluoranthene-d10	18.829	212	5441	0.166	ng	0.00
31) Terphenyl-d14	19.452	244	4123	0.266	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	13138	2.699	ng	98

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

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