

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034530.D
 Acq On : 10 Oct 2024 20:31
 Operator : RC/JU
 Sample : P4263-12
 Misc :
 ALS Vial : 106 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT162S1-20241001

Manual Integrations
 APPROVED

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

Quant Time: Oct 11 00:22:58 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	4173	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	11230	0.400	ng	0.01
13) Acenaphthene-d10	14.019	164	6541	0.400	ng	0.00
19) Phenanthrene-d10	16.784	188	12618	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	8198	0.400	ng	0.00
35) Perylene-d12	23.116	264	11902	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1371	0.116	ng	0.00
5) Phenol-d6	6.586	99	620m	0.045	ng	0.01
8) Nitrobenzene-d5	8.536	82	2229m	0.259	ng	0.02
11) 2-Methylnaphthalene-d10	11.738	152	5400	0.326	ng	0.01
14) 2,4,6-Tribromophenol	15.542	330	738	0.249	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	7948	0.299	ng	0.00
27) Fluoranthene-d10	18.829	212	13383	0.420	ng	0.00
31) Terphenyl-d14	19.447	244	9444	0.577	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.025	88	1259	0.241	ng	# 63

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

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