

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034540.D
 Acq On : 11 Oct 2024 03:50
 Operator : RC/JU
 Sample : P4241-03
 Misc :
 ALS Vial : 116 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT189D1-GW-20240927

Manual Integrations
 APPROVED

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

Quant Time: Oct 11 04:14:58 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	4059	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	10555	0.400	ng	0.01
13) Acenaphthene-d10	14.026	164	6182	0.400	ng	0.01
19) Phenanthrene-d10	16.791	188	11706	0.400	ng	0.01
29) Chrysene-d12	21.006	240	8588	0.400	ng	0.00
35) Perylene-d12	23.111	264	10546	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1203	0.104	ng	0.00
5) Phenol-d6	6.586	99	433m	0.032	ng	0.01
8) Nitrobenzene-d5	8.547	82	1923m	0.238	ng	0.03
11) 2-Methylnaphthalene-d10	11.746	152	4880m	0.313	ng	0.02
14) 2,4,6-Tribromophenol	15.549	330	642	0.229	ng	0.01
15) 2-Fluorobiphenyl	12.647	172	7365	0.293	ng	0.01
27) Fluoranthene-d10	18.832	212	11699	0.396	ng	0.00
31) Terphenyl-d14	19.445	244	11160	0.651	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.025	88	24863	4.900	ng	99
6) bis(2-Chloroethyl)ether	6.824	93	459	0.039	ng	93

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

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