

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034525.D
 Acq On : 10 Oct 2024 17:31
 Operator : RC/JU
 Sample : P4263-07
 Misc :
 ALS Vial : 101 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT188S1-20240930

Manual Integrations
 APPROVED

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

Quant Time: Oct 10 18:03:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 26 14:44:02 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	4207	0.400	ng	-0.02
7) Naphthalene-d8	10.127	136	11561	0.400	ng	#-0.02
13) Acenaphthene-d10	14.026	164	6606	0.400	ng	-0.02
19) Phenanthrene-d10	16.790	188	12259	0.400	ng	-0.01
29) Chrysene-d12	21.006	240	7970	0.400	ng	0.00
35) Perylene-d12	23.117	264	10230	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1917	0.161	ng	-0.01
5) Phenol-d6	6.593	99	1134m	0.082	ng	0.00
8) Nitrobenzene-d5	8.525	82	2408	0.272	ng	-0.01
11) 2-Methylnaphthalene-d10	11.731	152	5667	0.332	ng	-0.02
14) 2,4,6-Tribromophenol	15.537	330	879	0.294	ng	-0.01
15) 2-Fluorobiphenyl	12.636	172	8229	0.306	ng	-0.02
27) Fluoranthene-d10	18.832	212	12673	0.410	ng	-0.01
31) Terphenyl-d14	19.450	244	9755	0.613	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	730	0.139	ng	# 27

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

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