

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072724\
 Data File : BN033052.D
 Acq On : 28 Jul 2024 02:52
 Operator : MA/JU
 Sample : P3338-03
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 VPB196-HYD-20240718

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/29/2024
 Supervised By :mohammad ahmed 07/30/2024

Quant Time: Jul 29 01:11:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN071124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 21 23:33:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.561	152	2197	0.400	ng	# 0.03
7) Naphthalene-d8	10.372	136	9248m	0.400	ng	0.07
13) Acenaphthene-d10	14.190	164	6093	0.400	ng	0.01
19) Phenanthrene-d10	16.970	188	13064	0.400	ng	0.02
29) Chrysene-d12	21.169	240	11044	0.400	ng	0.00
35) Perylene-d12	23.367	264	9531	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.235	112	328m	0.059	ng	0.06
5) Phenol-d6	6.911	99	45m	0.007	ng	0.13
8) Nitrobenzene-d5	8.963	82	2893m	0.416	ng	0.23
11) 2-Methylnaphthalene-d10	12.007	152	11	0.001	ng	0.09
14) 2,4,6-Tribromophenol	15.753	330	640	0.246	ng	0.03
15) 2-Fluorobiphenyl	12.843	172	7718	0.340	ng	0.03
27) Fluoranthene-d10	18.997	212	7621	0.225	ng	0.00
31) Terphenyl-d14	19.582	244	10990	0.485	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.097	88	352	0.129	ng	# 53

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

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