

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072924\
 Data File : BN033065.D
 Acq On : 29 Jul 2024 12:04
 Operator : MA/JU
 Sample : P3338-03RE
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 VPB196-HYD-20240718RE

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/30/2024
 Supervised By :mohammad ahmed 07/30/2024

Quant Time: Jul 29 12:34:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN071124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 29 10:37:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.517	152	3358	0.400	ng	# 0.00
7) Naphthalene-d8	10.340	136	12141m	0.400	ng	0.01
13) Acenaphthene-d10	14.169	164	7621	0.400	ng	0.00
19) Phenanthrene-d10	16.945	188	17378	0.400	ng	#-0.01
29) Chrysene-d12	21.161	240	14614	0.400	ng	0.00
35) Perylene-d12	23.356	264	12550	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.221	112	523	0.062	ng	-0.02
5) Phenol-d6	6.903	99	71	0.007	ng	0.00
8) Nitrobenzene-d5	8.856	82	3657	0.400	ng	0.00
11) 2-Methylnaphthalene-d10	11.980	152	17	0.001	ng	0.03
14) 2,4,6-Tribromophenol	15.729	330	921	0.283	ng	-0.01
15) 2-Fluorobiphenyl	12.811	172	10376	0.366	ng	-0.01
27) Fluoranthene-d10	18.987	212	9870	0.220	ng	0.00
31) Terphenyl-d14	19.573	244	14375	0.480	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.097	88	378	0.091	ng	# 66

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

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