

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101124\
 Data File : BN034555.D
 Acq On : 11 Oct 2024 15:11
 Operator : RC/JU
 Sample : P4282-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW7-SP302-20241002

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/14/2024
 Supervised By :mohammad ahmed 10/15/2024

Quant Time: Oct 11 15:42:50 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	7.366	152	3890	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	10037	0.400	ng	0.01
13) Acenaphthene-d10	14.019	164	5820	0.400	ng	0.00
19) Phenanthrene-d10	16.784	188	10479	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	6403	0.400	ng	0.00
35) Perylene-d12	23.110	264	9770	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	1362	0.123	ng	0.00
5) Phenol-d6	6.586	99	595m	0.046	ng	0.01
8) Nitrobenzene-d5	8.536	82	2406m	0.313	ng	0.02
11) 2-Methylnaphthalene-d10	11.742	152	5011	0.338	ng	0.02
14) 2,4,6-Tribromophenol	15.555	330	352	0.133	ng	0.02
15) 2-Fluorobiphenyl	12.640	172	7599	0.321	ng	0.00
27) Fluoranthene-d10	18.825	212	11831	0.447	ng	0.00
31) Terphenyl-d14	19.443	244	9331	0.730	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	355	0.073	ng	# 36

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

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