

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101124\
 Data File : BN034557.D
 Acq On : 11 Oct 2024 16:23
 Operator : RC/JU
 Sample : P4282-01DL 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW7-SP100-20241002DL

Manual Integrations
 APPROVED

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

Quant Time: Oct 11 17:27:47 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	7.365	152	3752	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	10492	0.400	ng	0.01
13) Acenaphthene-d10	14.019	164	6659	0.400	ng	0.00
19) Phenanthrene-d10	16.783	188	12028	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	6979	0.400	ng	0.00
35) Perylene-d12	23.110	264	10316	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	645	0.061	ng	0.00
5) Phenol-d6	6.650	99	296m	0.024	ng	0.08
8) Nitrobenzene-d5	8.557	82	935m	0.117	ng	0.04
11) 2-Methylnaphthalene-d10	11.765	152	2266m	0.146	ng	0.04
14) 2,4,6-Tribromophenol	15.555	330	298	0.099	ng	0.02
15) 2-Fluorobiphenyl	12.651	172	3395	0.125	ng	0.01
27) Fluoranthene-d10	18.825	212	6100	0.201	ng	0.00
31) Terphenyl-d14	19.443	244	5726	0.411	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.018	88	13245	2.824	ng	99

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

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