

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034537.D
 Acq On : 11 Oct 2024 02:02
 Operator : RC/JU
 Sample : PB163903BL
 Misc :
 ALS Vial : 113 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB163903BL

Manual Integrations
 APPROVED

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

Quant Time: Oct 11 02:31:14 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	7.380	152	3859	0.400	ng	0.00
7) Naphthalene-d8	10.148	136	10903	0.400	ng	0.02
13) Acenaphthene-d10	14.030	164	6019	0.400	ng	0.01
19) Phenanthrene-d10	16.808	188	10284	0.400	ng	# 0.03
29) Chrysene-d12	21.017	240	6672	0.400	ng	0.01
35) Perylene-d12	23.119	264	8124	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	3793	0.346	ng	0.00
5) Phenol-d6	6.614	99	3439m	0.270	ng	0.04
8) Nitrobenzene-d5	8.557	82	2354m	0.282	ng	0.04
11) 2-Methylnaphthalene-d10	11.761	152	5813m	0.361	ng	0.03
14) 2,4,6-Tribromophenol	15.567	330	638m	0.234	ng	0.03
15) 2-Fluorobiphenyl	12.661	172	8252m	0.337	ng	0.03
27) Fluoranthene-d10	18.839	212	9697	0.374	ng	0.01
31) Terphenyl-d14	19.456	244	6149	0.461	ng	0.00

Target Compounds Qvalue

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

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