

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101124\
 Data File : BN034553.D
 Acq On : 11 Oct 2024 13:59
 Operator : RC/JU
 Sample : PB163903BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB163903BL

Manual Integrations
 APPROVED

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

Quant Time: Oct 11 15:42:12 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	4285	0.400	ng	0.00
7) Naphthalene-d8	10.127	136	11473	0.400	ng	# 0.00
13) Acenaphthene-d10	14.015	164	6303	0.400	ng	0.00
19) Phenanthrene-d10	16.790	188	10680	0.400	ng	0.01
29) Chrysene-d12	21.006	240	6687	0.400	ng	0.00
35) Perylene-d12	23.111	264	10014	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	4065	0.334	ng	0.00
5) Phenol-d6	6.585	99	3635m	0.257	ng	0.01
8) Nitrobenzene-d5	8.536	82	2636m	0.300	ng	0.02
11) 2-Methylnaphthalene-d10	11.738	152	5997m	0.354	ng	0.01
14) 2,4,6-Tribromophenol	15.549	330	396	0.139	ng	0.01
15) 2-Fluorobiphenyl	12.647	172	7594	0.296	ng	0.01
27) Fluoranthene-d10	18.827	212	10366	0.385	ng	0.00
31) Terphenyl-d14	19.445	244	6932	0.519	ng	0.00

Target Compounds Qvalue

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

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