

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070924\
 Data File : BN032565.D
 Acq On : 09 Jul 2024 00:39
 Operator : MA/JU
 Sample : PB161896BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB161896BL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/09/2024
 Supervised By :mohammad ahmed 07/11/2024

Quant Time: Jul 09 03:13:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN070924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 09 02:39:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.610	152	3500m	0.400	ng	0.02
7) Naphthalene-d8	10.442	136	11577m	0.400	ng	0.06
13) Acenaphthene-d10	14.263	164	6937	0.400	ng	0.03
19) Phenanthrene-d10	17.053	188	14116	0.400	ng	# 0.05
29) Chrysene-d12	21.238	240	13062	0.400	ng	# 0.03
35) Perylene-d12	23.443	264	12388	0.400	ng	# 0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.219	112	3789	0.428	ng	0.00
5) Phenol-d6	6.924	99	4254m	0.376	ng	0.10
8) Nitrobenzene-d5	8.926	82	3336m	0.386	ng	0.13
11) 2-Methylnaphthalene-d10	12.069	152	7083m	0.387	ng	0.09
14) 2,4,6-Tribromophenol	15.825	330	573m	0.182	ng	0.06
15) 2-Fluorobiphenyl	12.948	172	10593m	0.414	ng	0.09
27) Fluoranthene-d10	19.063	212	17059	0.464	ng	0.02
31) Terphenyl-d14	19.662	244	12234	0.384	ng	0.01

Target Compounds Qvalue

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

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