

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072724\
 Data File : BN033056.D
 Acq On : 28 Jul 2024 06:00
 Operator : MA/JU
 Sample : PB162236BL
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB162236BL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/29/2024
 Supervised By :mohammad ahmed 07/30/2024

Quant Time: Jul 29 01:12:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN071124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 21 23:33:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.589	152	2873m	0.400	ng	0.06
7) Naphthalene-d8	10.490	136	9361m	0.400	ng	0.19
13) Acenaphthene-d10	14.229	164	5708m	0.400	ng	0.05
19) Phenanthrene-d10	17.026	188	10377m	0.400	ng	0.07
29) Chrysene-d12	21.194	240	9562	0.400	ng	# 0.03
35) Perylene-d12	23.377	264	9511	0.400	ng	# 0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.228	112	2987m	0.411	ng	0.05
5) Phenol-d6	7.120	99	3670m	0.407	ng	0.34
8) Nitrobenzene-d5	9.059	82	2898m	0.411	ng	0.32
11) 2-Methylnaphthalene-d10	12.094	152	5012m	0.350	ng	0.18
14) 2,4,6-Tribromophenol	15.822	330	513m	0.210	ng	0.10
15) 2-Fluorobiphenyl	12.946	172	7434m	0.350	ng	0.14
27) Fluoranthene-d10	19.031	212	11518	0.429	ng	0.04
31) Terphenyl-d14	19.612	244	8508	0.434	ng	0.02

Target Compounds Qvalue

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

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