

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034138.D
 Acq On : 21 Sep 2024 23:58
 Operator : JU/RC
 Sample : PB163558BL
 Misc :
 ALS Vial : 103 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB163558BL

Quant Time: Sep 23 00:01:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.430	152	5847	0.400	ng	0.00
7) Naphthalene-d8	10.180	136	15255	0.400	ng	#-0.02
13) Acenaphthene-d10	14.072	164	8244	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	15908	0.400	ng	# 0.00
29) Chrysene-d12	21.053	240	8486	0.400	ng	0.00
35) Perylene-d12	23.174	264	7265	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	5334	0.321	ng	0.00
5) Phenol-d6	6.629	99	5074	0.263	ng	0.00
8) Nitrobenzene-d5	8.568	82	3655	0.313	ng	0.00
11) 2-Methylnaphthalene-d10	11.787	152	7228	0.321	ng	0.00
14) 2,4,6-Tribromophenol	15.592	330	498	0.133	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	11070	0.330	ng	0.00
27) Fluoranthene-d10	18.876	212	13278	0.331	ng	-0.01
31) Terphenyl-d14	19.494	244	7712	0.455	ng	-0.01

Target Compounds Qvalue

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

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