

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081724\
 Data File : BN033429.D
 Acq On : 16 Aug 2024 14:10
 Operator : MA/JU
 Sample : PB162692BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB162692BL

Quant Time: Aug 16 14:47:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 16 04:44:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.567	152	7905	0.400	ng	0.00
7) Naphthalene-d8	10.325	136	21401	0.400	ng	-0.01
13) Acenaphthene-d10	14.199	164	10134	0.400	ng	0.00
19) Phenanthrene-d10	16.942	188	20177	0.400	ng	#-0.01
29) Chrysene-d12	21.157	240	15595	0.400	ng	0.00
35) Perylene-d12	23.332	264	17788	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.198	112	8291	0.380	ng	0.00
5) Phenol-d6	6.750	99	9111	0.309	ng	0.00
8) Nitrobenzene-d5	8.702	82	6945	0.430	ng	0.00
11) 2-Methylnaphthalene-d10	11.926	152	11257	0.342	ng	0.00
14) 2,4,6-Tribromophenol	15.700	330	1085	0.209	ng	0.00
15) 2-Fluorobiphenyl	12.820	172	14855	0.358	ng	0.00
27) Fluoranthene-d10	18.989	212	18085	0.335	ng	0.00
31) Terphenyl-d14	19.607	244	13399	0.455	ng	0.00

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

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