

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110122\
 Data File : BN022446.D
 Acq On : 01 Nov 2022 23:35
 Operator : CG/JU
 Sample : PB148793BL
 Misc :
 ALS Vial : 54 Sample Multiplier: 50

Instrument :
 BNA_N
 ClientSampleId :
 PB148793BL

Quant Time: Nov 02 06:04:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 02 05:31:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.921	152	175905	20.000	ng	0.00
7) Naphthalene-d8	10.752	136	591852	20.000	ng	# 0.00
13) Acenaphthene-d10	14.589	164	307968	20.000	ng	0.00
19) Phenanthrene-d10	17.350	188	533506	20.000	ng	# 0.00
29) Chrysene-d12	21.546	240	300087	20.000	ng	0.00
35) Perylene-d12	23.996	264	260470	20.000	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.458	112	1544462	164.746	ng	0.00
5) Phenol-d6	7.090	99	2095201	178.367	ng	0.00
8) Nitrobenzene-d5	9.097	82	1051274	102.696	ng	-0.01
11) 2-Methylnaphthalene-d10	12.357	152	21	0.001	ng	0.00
14) 2,4,6-Tribromophenol	16.096	330	473399	178.833	ng	0.00
15) 2-Fluorobiphenyl	13.221	172	2151269	81.659	ng	0.00
27) Fluoranthene-d10	19.395	212	5	0.000	ng	0.00
31) Terphenyl-d14	19.985	244	1686145	107.851	ng	0.00

Target Compounds Qvalue

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

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