

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102224\
 Data File : BN034682.D
 Acq On : 22 Oct 2024 18:29
 Operator : RC/JU
 Sample : PB164282BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB164282BL

Quant Time: Oct 23 00:48:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 22 16:43:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.662	152	9218	0.400	ng	0.00
7) Naphthalene-d8	10.436	136	25811	0.400	ng	0.00
13) Acenaphthene-d10	14.293	164	12398	0.400	ng	0.00
19) Phenanthrene-d10	17.039	188	25006	0.400	ng	# 0.00
29) Chrysene-d12	21.221	240	15417	0.400	ng	0.00
35) Perylene-d12	23.426	264	5086	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.278	112	8542	0.300	ng	0.00
5) Phenol-d6	6.838	99	10532	0.291	ng	0.00
8) Nitrobenzene-d5	8.803	82	7168	0.334	ng	0.00
11) 2-Methylnaphthalene-d10	12.030	152	12261	0.334	ng	0.00
14) 2,4,6-Tribromophenol	15.785	330	1316	0.240	ng	0.00
15) 2-Fluorobiphenyl	12.914	172	16781	0.342	ng	0.00
27) Fluoranthene-d10	19.064	212	20381	0.326	ng	0.00
31) Terphenyl-d14	19.677	244	12380	0.417	ng	0.00

Target Compounds Qvalue

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

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