

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071025\
 Data File : BN037463.D
 Acq On : 10 Jul 2025 19:26
 Operator : RC/JU
 Sample : PB168696BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB168696BL

Quant Time: Jul 11 01:28:56 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 10 00:48:38 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.732	152	1779	0.400	ng	0.00
7) Naphthalene-d8	10.509	136	4164	0.400	ng	-0.01
13) Acenaphthene-d10	14.366	164	1917	0.400	ng	0.00
19) Phenanthrene-d10	17.099	188	3331	0.400	ng	-0.01
29) Chrysene-d12	21.286	240	2383	0.400	ng	# 0.00
35) Perylene-d12	23.528	264	2255	0.400	ng	#-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	1492	0.312	ng	0.00
5) Phenol-d6	6.887	99	1835	0.311	ng	0.00
8) Nitrobenzene-d5	8.865	82	1171	0.418	ng	-0.01
11) 2-Methylnaphthalene-d10	12.106	152	2033	0.349	ng	0.00
14) 2,4,6-Tribromophenol	15.845	330	118	0.151	ng	-0.01
15) 2-Fluorobiphenyl	12.988	172	4155	0.393	ng	0.00
27) Fluoranthene-d10	19.132	212	2941	0.343	ng	0.00
31) Terphenyl-d14	19.740	244	2014	0.401	ng	0.00

Target Compounds Qvalue

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

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