

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062725\
 Data File : BN037418.D
 Acq On : 27 Jun 2025 10:45
 Operator : RC/JU
 Sample : PB168563BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB168563BL

Quant Time: Jun 27 11:14:05 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN062625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 26 16:06:33 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.568	152	1980	0.400	ng	0.00
7) Naphthalene-d8	10.340	136	3978	0.400	ng	# 0.00
13) Acenaphthene-d10	14.213	164	2680	0.400	ng	0.00
19) Phenanthrene-d10	16.971	188	5035	0.400	ng	0.01
29) Chrysene-d12	21.171	240	4921	0.400	ng	0.00
35) Perylene-d12	23.348	264	3783	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.163	112	1439	0.377	ng	0.00
5) Phenol-d6	6.751	99	1135	0.287	ng	0.00
8) Nitrobenzene-d5	8.728	82	1008	0.316	ng	0.01
11) 2-Methylnaphthalene-d10	11.945	152	2119	0.344	ng	0.00
14) 2,4,6-Tribromophenol	15.730	330	523	0.333	ng	0.01
15) 2-Fluorobiphenyl	12.838	172	4000	0.346	ng	0.00
27) Fluoranthene-d10	19.003	212	5607	0.389	ng	0.00
31) Terphenyl-d14	19.611	244	4125	0.393	ng	0.00

Target Compounds Qvalue

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

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