

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091725\
 Data File : BN037741.D
 Acq On : 17 Sep 2025 12:02
 Operator : RC/JU
 Sample : PB169692BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB169692BL

Quant Time: Sep 17 12:51:54 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 11 14:20:01 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	7434	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	20202	0.400	ng	# 0.00
13) Acenaphthene-d10	14.494	164	8991	0.400	ng	0.00
19) Phenanthrene-d10	17.235	188	14614	0.400	ng	0.00
29) Chrysene-d12	21.411	240	12468	0.400	ng	# 0.00
35) Perylene-d12	23.750	264	13022	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	5588	0.327	ng	0.00
5) Phenol-d6	6.988	99	6691	0.302	ng	0.00
8) Nitrobenzene-d5	8.982	82	4866	0.329	ng	-0.01
11) 2-Methylnaphthalene-d10	12.238	152	8483	0.317	ng	0.00
14) 2,4,6-Tribromophenol	15.982	330	473	0.202	ng	0.00
15) 2-Fluorobiphenyl	13.115	172	13040	0.347	ng	0.00
27) Fluoranthene-d10	19.271	212	10845	0.296	ng	0.00
31) Terphenyl-d14	19.870	244	8091	0.320	ng	0.00

Target Compounds	Qvalue

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

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