

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110725\
 Data File : BN038166.D
 Acq On : 07 Nov 2025 22:54
 Operator : RC/JU
 Sample : PB170427BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB170427BL

Quant Time: Nov 08 03:28:02 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 29 13:09:46 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	8786	0.400	ng	0.00
7) Naphthalene-d8	10.573	136	25751	0.400	ng	# 0.00
13) Acenaphthene-d10	14.441	164	13576	0.400	ng	0.01
19) Phenanthrene-d10	17.186	188	20711	0.400	ng	# 0.00
29) Chrysene-d12	21.375	240	12967	0.400	ng	# 0.00
35) Perylene-d12	23.686	264	14107	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.348	112	8397	0.406	ng	0.00
5) Phenol-d6	6.937	99	9322	0.370	ng	0.00
8) Nitrobenzene-d5	8.929	82	8179	0.394	ng	0.00
11) 2-Methylnaphthalene-d10	12.177	152	13506	0.367	ng	0.00
14) 2,4,6-Tribromophenol	15.945	330	1265	0.226	ng	0.01
15) 2-Fluorobiphenyl	13.062	172	22627	0.416	ng	0.01
27) Fluoranthene-d10	19.225	212	16467	0.315	ng	0.00
31) Terphenyl-d14	19.824	244	12995	0.461	ng	0.00

Target Compounds Qvalue

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

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