

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092425\
 Data File : BN037877.D
 Acq On : 24 Sep 2025 15:48
 Operator : RC/JU
 Sample : Q3015-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 WP0925-PT-ACIDS-WP

Quant Time: Sep 24 16:19:06 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN092325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 24 11:00:01 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.833	152	5760	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	17583	0.400	ng	0.01
13) Acenaphthene-d10	14.484	164	8737	0.400	ng	0.01
19) Phenanthrene-d10	17.223	188	16666	0.400	ng	# 0.00
29) Chrysene-d12	21.411	240	12568	0.400	ng	0.00
35) Perylene-d12	23.741	264	11362	0.400	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	2442475	185.820	ng	0.01
5) Phenol-d6	6.995	99	3401287	196.945	ng	0.02
8) Nitrobenzene-d5	8.982	82	1681652	125.388	ng	0.01
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng	
14) 2,4,6-Tribromophenol	15.969	330	1057879	319.317	ng	0.00
15) 2-Fluorobiphenyl	13.105	172	4153453	116.120	ng	0.01
27) Fluoranthene-d10	0.000	212	0d	0.000	ng	
31) Terphenyl-d14	19.861	244	3706695	148.066	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.597	198	398438	33.153	ng	# 86
24) Pentachlorophenol	16.875	266	537511	118.961	ng	97

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

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