

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092425\
 Data File : BN037878.D
 Acq On : 24 Sep 2025 16:58
 Operator : RC/JU
 Sample : Q3015-08DL 50X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Quant Time: Sep 24 17:24:45 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	5448	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	16730	0.400	ng	0.01
13) Acenaphthene-d10	14.484	164	8546	0.400	ng	0.01
19) Phenanthrene-d10	17.223	188	17144	0.400	ng	# 0.00
29) Chrysene-d12	21.411	240	13570	0.400	ng	0.00
35) Perylene-d12	23.741	264	12821	0.400	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	41384	3.329	ng	0.01
5) Phenol-d6	6.988	99	55546	3.400	ng	0.01
8) Nitrobenzene-d5	8.982	82	29544	2.315	ng	0.01
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng	
14) 2,4,6-Tribromophenol	15.969	330	11197	3.455	ng	0.00
15) 2-Fluorobiphenyl	13.105	172	85599	2.447	ng	0.01
27) Fluoranthene-d10	19.262	212	8	0.000	ng	0.00
31) Terphenyl-d14	19.861	244	76232	2.820	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.597	198	2740	1.185	ng	# 88
24) Pentachlorophenol	16.875	266	4892	1.053	ng	97

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

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