

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031425\
 Data File : BN036608.D
 Acq On : 14 Mar 2025 15:33
 Operator : RC/JU
 Sample : Q1502-08DL 20X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 14 16:00:00 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 16:06:28 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.717	152	2324	0.400	ng	0.00
7) Naphthalene-d8	10.498	136	6014	0.400	ng	-0.01
13) Acenaphthene-d10	14.366	164	3371	0.400	ng	0.00
19) Phenanthrene-d10	17.111	188	7010	0.400	ng	0.00
29) Chrysene-d12	21.304	240	4582	0.400	ng	0.00
35) Perylene-d12	23.560	264	3725	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.305	112	41988	7.752	ng	0.00
5) Phenol-d6	6.887	99	54884	8.204	ng	-0.01
8) Nitrobenzene-d5	8.865	82	30840	4.714	ng	-0.01
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng	
14) 2,4,6-Tribromophenol	15.845	330	13092	8.559	ng	-0.01
15) 2-Fluorobiphenyl	12.983	172	105143	5.362	ng	0.00
27) Fluoranthene-d10	19.127	212	2	0.000	ng	-0.01
31) Terphenyl-d14	19.740	244	68962	6.282	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.489	198	8530	4.417	ng	Qvalue # 46
24) Pentachlorophenol	16.764	266	10797	4.465	ng	99

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

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