

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081325\
 Data File : BP025387.D
 Acq On : 13 Aug 2025 16:50
 Operator : CG/JU
 Sample : Q2820-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 82H-N

Quant Time: Aug 13 17:11:54 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	177935	20.000	ng	0.00
21) Naphthalene-d8	10.566	136	679453	20.000	ng	0.00
39) Acenaphthene-d10	14.413	164	421690	20.000	ng	0.00
64) Phenanthrene-d10	17.207	188	844698	20.000	ng	-0.02
76) Chrysene-d12	21.666	240	1026814	20.000	ng	0.00
86) Perylene-d12	25.060	264	1145461	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.408	112	815476	73.711	ng	0.00
7) Phenol-d6	6.966	99	973941	66.892	ng	0.00
23) Nitrobenzene-d5	8.931	82	642226	52.366	ng	-0.01
42) 2,4,6-Tribromophenol	15.925	330	424930	101.009	ng	0.00
45) 2-Fluorobiphenyl	13.025	172	1499283	49.108	ng	0.00
79) Terphenyl-d14	19.930	244	2514493	46.253	ng	-0.01
Target Compounds						
75) Fluoranthene	19.330	202	155534	2.904	ng	98
78) Pyrene	19.707	202	156861	2.297	ng	98

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

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