

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062725\
 Data File : BF142901.D
 Acq On : 27 Jun 2025 19:10
 Operator : RC/JU
 Sample : Q2420-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11933

Quant Time: Jun 27 22:40:35 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 24 05:28:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	56236	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	196491	20.000	ng	-0.01
39) Acenaphthene-d10	9.916	164	89932	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	147476	20.000	ng	0.00
76) Chrysene-d12	14.051	240	149469	20.000	ng	0.00
86) Perylene-d12	15.545	264	103761	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	386448	114.412	ng	0.01
7) Phenol-d6	6.510	99	390633	98.025	ng	0.00
23) Nitrobenzene-d5	7.439	82	290911	81.208	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	121600	131.394	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	592366	86.529	ng	-0.01
79) Terphenyl-d14	12.992	244	649160	60.009	ng	0.00
Target Compounds						
						Qvalue
52) Acenaphthene	9.951	154	21778	4.067	ng	99
71) Phenanthrene	11.427	178	37350	4.675	ng	98
75) Fluoranthene	12.621	202	22066	2.910	ng	99

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

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