

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031125\
 Data File : BF141923.D
 Acq On : 11 Mar 2025 18:42
 Operator : RC/JU
 Sample : Q1523-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-A1-01-C

Quant Time: Mar 12 01:15:12 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 15:46:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	132841	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	498756	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	254340	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	415187	20.000	ng	0.00
76) Chrysene-d12	14.033	240	364222	20.000	ng	0.00
86) Perylene-d12	15.504	264	275509	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	988695	124.216	ng	0.01
7) Phenol-d6	6.504	99	1099758	108.519	ng	0.00
23) Nitrobenzene-d5	7.440	82	844353	95.271	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	436097	135.155	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	1723615	103.063	ng	0.00
79) Terphenyl-d14	12.986	244	2030864	82.437	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.181	128	545492	21.291	ng	100
37) 2-Methylnaphthalene	8.869	142	103047	6.017	ng	98
38) 1-Methylnaphthalene	8.969	142	125009	7.565	ng	99
52) Acenaphthene	9.945	154	35731	2.379	ng	97
77) Benzidine	12.733	184	77233	15.199	ng	98

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

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