

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052225\
 Data File : BP024779.D
 Acq On : 23 May 2025 05:15
 Operator : RC/JU
 Sample : Q2079-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-A2-03-C

Quant Time: May 23 05:49:23 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 13 16:21:39 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	90311	20.000	ng	-0.01
21) Naphthalene-d8	10.422	136	341108	20.000	ng	#-0.01
39) Acenaphthene-d10	14.281	164	209629	20.000	ng	-0.02
64) Phenanthrene-d10	17.081	188	421419	20.000	ng	-0.02
76) Chrysene-d12	21.516	240	510159	20.000	ng	0.00
86) Perylene-d12	24.810	264	663924	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.287	112	601055	113.832	ng	0.00
7) Phenol-d6	6.852	99	640185	94.997	ng	0.00
23) Nitrobenzene-d5	8.805	82	469306	69.516	ng	0.00
42) 2,4,6-Tribromophenol	15.804	330	530300	149.220	ng	0.00
45) 2-Fluorobiphenyl	12.893	172	1241727	77.605	ng	-0.01
79) Terphenyl-d14	19.804	244	2387814	80.257	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.475	128	6746808	381.682	ng	99
37) 2-Methylnaphthalene	12.081	142	776000	67.798	ng	99
38) 1-Methylnaphthalene	12.304	142	645720	52.726	ng	100
46) 1,1'-Biphenyl	13.104	154	118692	7.661	ng	99
52) Acenaphthene	14.346	154	339474	29.869	ng	100
58) Fluorene	15.351	166	125768	8.508	ng	100
71) Phenanthrene	17.122	178	243735	10.404	ng	99

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

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