

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040422\
 Data File : BP009686.D
 Acq On : 04 Apr 2022 19:15
 Operator : CG/JU
 Sample : N2227-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 124055

Quant Time: Apr 04 23:27:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 31 04:36:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	250732	20.000	ng	-0.01
21) Naphthalene-d8	10.516	136	996671	20.000	ng	-0.02
39) Acenaphthene-d10	14.375	164	621347	20.000	ng	-0.02
64) Phenanthrene-d10	17.139	188	1291194	20.000	ng	-0.02
76) Chrysene-d12	21.263	240	1294336	20.000	ng	-0.01
86) Perylene-d12	23.633	264	1356823	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	2021744	140.781	ng	-0.01
7) Phenol-d6	6.922	99	2538750	130.721	ng	-0.01
23) Nitrobenzene-d5	8.881	82	1617189	86.225	ng	-0.02
42) 2,4,6-Tribromophenol	15.881	330	1072459	104.596	ng	-0.01
45) 2-Fluorobiphenyl	12.993	172	3428888	76.502	ng	-0.02
79) Terphenyl-d14	19.722	244	5413911	75.076	ng	-0.02
Target Compounds						
						Qvalue
75) Fluoranthene	19.169	202	360653	4.310	ng	99
78) Pyrene	19.522	202	495191	5.806	ng	100
81) Benzo(a)anthracene	21.245	228	264700	3.113	ng	99
83) Chrysene	21.298	228	254589	3.162	ng	98
88) Benzo(b)fluoranthene	22.916	252	350431	4.326	ng	99
90) Benzo(a)pyrene	23.527	252	169764	2.450	ng	99

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

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