

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040422\
 Data File : BP009687.D
 Acq On : 04 Apr 2022 19:49
 Operator : CG/JU
 Sample : N2227-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 124035

Quant Time: Apr 04 23:27:17 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	7.722	152	252656	20.000	ng	-0.02
21) Naphthalene-d8	10.516	136	999543	20.000	ng	-0.02
39) Acenaphthene-d10	14.375	164	633821	20.000	ng	-0.02
64) Phenanthrene-d10	17.139	188	1299385	20.000	ng	-0.02
76) Chrysene-d12	21.257	240	1315898	20.000	ng	-0.02
86) Perylene-d12	23.633	264	1373303	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	1869554	129.192	ng	-0.01
7) Phenol-d6	6.922	99	2321496	118.625	ng	-0.01
23) Nitrobenzene-d5	8.881	82	1465299	77.902	ng	-0.02
42) 2,4,6-Tribromophenol	15.881	330	971525	92.888	ng	-0.01
45) 2-Fluorobiphenyl	12.993	172	3121717	68.278	ng	-0.02
79) Terphenyl-d14	19.722	244	4994872	68.130	ng	-0.02
Target Compounds						
75) Fluoranthene	19.169	202	206751	2.455	ng	Qvalue 99

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

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