

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
 Data File : BP009427.D
 Acq On : 16 Mar 2022 17:46
 Operator : CG/JU
 Sample : N1952-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CON-1

Quant Time: Mar 17 01:09:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	393393	20.000	ng	-0.01
21) Naphthalene-d8	10.605	136	1526429	20.000	ng	-0.01
39) Acenaphthene-d10	14.451	164	900207	20.000	ng	-0.01
64) Phenanthrene-d10	17.210	188	1513516	20.000	ng	-0.01
76) Chrysene-d12	21.322	240	1249212	20.000	ng	0.00
86) Perylene-d12	23.733	264	1557645	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	2661683	98.762	ng	0.00
7) Phenol-d6	7.005	99	3454673	100.871	ng	0.02
23) Nitrobenzene-d5	8.964	82	2280392	82.284	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	1331370	122.481	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	5425460	82.882	ng	-0.01
79) Terphenyl-d14	19.786	244	5505711	83.172	ng	-0.01
Target Compounds						
						Qvalue
60) Diethylphthalate	15.281	149	192585	2.979	ng	98
78) Pyrene	19.586	202	319353	3.786	ng	98
83) Chrysene	21.357	228	173525	2.111	ng	# 90

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

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