

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032922\
 Data File : BP009604.D
 Acq On : 29 Mar 2022 20:45
 Operator : CG/JU
 Sample : N1949-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CP-BH2-031422-4-6

Quant Time: Mar 30 04:10:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 30 04:00:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	280036	20.000	ng	0.00
21) Naphthalene-d8	10.552	136	1122817	20.000	ng	0.00
39) Acenaphthene-d10	14.404	164	707668	20.000	ng	0.00
64) Phenanthrene-d10	17.169	188	1488754	20.000	ng	0.00
76) Chrysene-d12	21.286	240	1630629	20.000	ng	0.00
86) Perylene-d12	23.680	264	1788380	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.364	112	1491301	92.978	ng	0.00
7) Phenol-d6	6.952	99	1938080	89.350	ng	0.00
23) Nitrobenzene-d5	8.916	82	1181676	55.926	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	856783	73.369	ng	0.00
45) 2-Fluorobiphenyl	13.028	172	2841262	55.659	ng	0.00
79) Terphenyl-d14	19.751	244	4863878	53.538	ng	0.00
Target Compounds						
84) Bis(2-ethylhexyl)phtha...	21.204	149	174331	2.739	ng	Qvalue 100
88) Benzo(b)fluoranthene	22.951	252	270408	2.533	ng	# 85
89) Benzo(k)fluoranthene	23.004	252	283373	2.711	ng	# 91
90) Benzo(a)pyrene	23.574	252	210772	2.307	ng	# 90

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

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