

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060222\
 Data File : BF128881.D
 Acq On : 20 Jun 2022 17:21
 Operator : CG\JU
 Sample : N3373-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-1S

Quant Time: Jun 21 01:16:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 21 00:40:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	109898	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	420707	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	199841	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	363385	20.000	ng	0.00
76) Chrysene-d12	13.945	240	218346	20.000	ng	# 0.00
86) Perylene-d12	15.392	264	172670	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	599428	86.879	ng	0.00
7) Phenol-d6	6.428	99	921328	109.394	ng	-0.01
23) Nitrobenzene-d5	7.339	82	673817	74.570	ng	-0.01
42) 2,4,6-Tribromophenol	10.604	330	81780	34.635	ng	0.00
45) 2-Fluorobiphenyl	9.133	172	1119336	78.461	ng	0.00
79) Terphenyl-d14	12.892	244	1177536	93.710	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.322	178	381070	20.907	ng	98
73) Carbazole	11.533	167	52795	3.151	ng	99
75) Fluoranthene	12.522	202	1188347	62.373	ng	99
78) Pyrene	12.745	202	187272	11.528	ng	96

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

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