

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062022\
 Data File : BF128889.D
 Acq On : 20 Jun 2022 21:14
 Operator : CG\JU
 Sample : N3373-02DL 20X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-1EDL

Quant Time: Jun 21 01:20:55 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	129252	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	505771	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	290907	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	542825	20.000	ng	0.00
76) Chrysene-d12	13.939	240	322320	20.000	ng	#-0.01
86) Perylene-d12	15.392	264	214900	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	42497	5.237	ng	0.00
7) Phenol-d6	6.428	99	56934	5.748	ng	-0.01
23) Nitrobenzene-d5	7.334	82	40339	3.713	ng	-0.02
42) 2,4,6-Tribromophenol	10.598	330	11898	3.462	ng	-0.01
45) 2-Fluorobiphenyl	9.128	172	88310	4.252	ng	-0.01
79) Terphenyl-d14	12.880	244	96288	5.191	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.322	178	464438	17.058	ng	100
73) Carbazole	11.533	167	85197	3.404	ng	97
75) Fluoranthene	12.522	202	1389132	48.809	ng	95
78) Pyrene	12.745	202	736216	30.701	ng	99

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

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