

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060222\
 Data File : BF128886.D
 Acq On : 20 Jun 2022 19:47
 Operator : CG\JU
 Sample : N3373-05DL 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-6SDL

Quant Time: Jun 21 01:19:37 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	131413	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	424383	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	253000	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	466057	20.000	ng	0.00
76) Chrysene-d12	13.939	240	297237	20.000	ng	#-0.01
86) Perylene-d12	15.392	264	209125	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	137090	16.616	ng	0.00
7) Phenol-d6	6.428	99	190446	18.911	ng	-0.01
23) Nitrobenzene-d5	7.333	82	123959	13.599	ng	-0.02
42) 2,4,6-Tribromophenol	10.604	330	13164	4.404	ng	0.00
45) 2-Fluorobiphenyl	9.127	172	243317	13.472	ng	-0.01
79) Terphenyl-d14	12.880	244	310984	18.180	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.327	178	889686	38.059	ng	99
73) Carbazole	11.533	167	51980	2.419	ng	99
75) Fluoranthene	12.521	202	1639445	67.093	ng	98
78) Pyrene	12.745	202	830495	37.555	ng	100

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

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