

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060222\
 Data File : BF128887.D
 Acq On : 20 Jun 2022 20:16
 Operator : CG\JU
 Sample : N3373-03DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-1WDL

Quant Time: Jun 21 01:20:06 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	133301	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	522424	20.000	ng	0.00
39) Acenaphthene-d10	9.810	164	303270	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	545675	20.000	ng	0.00
76) Chrysene-d12	13.939	240	313917	20.000	ng	#-0.01
86) Perylene-d12	15.392	264	249369	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	126081	15.066	ng	0.00
7) Phenol-d6	6.422	99	179863	17.607	ng	-0.02
23) Nitrobenzene-d5	7.334	82	127741	11.384	ng	-0.02
42) 2,4,6-Tribromophenol	10.604	330	9077	2.533	ng	0.00
45) 2-Fluorobiphenyl	9.128	172	267448	12.353	ng	-0.01
79) Terphenyl-d14	12.880	244	280275	15.514	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.322	178	420019	15.346	ng	99
73) Carbazole	11.533	167	131312	5.219	ng	97
75) Fluoranthene	12.521	202	1721028	60.155	ng	98
78) Pyrene	12.745	202	898095	38.454	ng	99

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

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