

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
 Data File : BF128912.D
 Acq On : 21 Jun 2022 22:24
 Operator : CG\JU
 Sample : N3373-08DL 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-6WDL

Quant Time: Jun 22 01:56:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 01:02:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.751	152	108593	20.000	ng	0.00
21) Naphthalene-d8	8.039	136	390316	20.000	ng	# 0.00
39) Acenaphthene-d10	9.792	164	186239	20.000	ng	0.00
64) Phenanthrene-d10	11.274	188	281738	20.000	ng	0.00
76) Chrysene-d12	13.921	240	239979	20.000	ng	# 0.00
86) Perylene-d12	15.362	264	168344	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.392	112	50245	7.370	ng	0.00
7) Phenol-d6	6.416	99	62363	7.494	ng	0.00
23) Nitrobenzene-d5	7.316	82	45326	5.407	ng	-0.01
42) 2,4,6-Tribromophenol	10.580	330	10099	4.589	ng	0.00
45) 2-Fluorobiphenyl	9.110	172	85845	6.457	ng	0.00
79) Terphenyl-d14	12.863	244	73083	5.292	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.298	178	201456	14.256	ng	98
73) Carbazole	11.510	167	36992	2.847	ng	99
75) Fluoranthene	12.498	202	624925	42.306	ng	94
78) Pyrene	12.721	202	372257	20.850	ng	98

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

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