

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062022\
 Data File : BF128888.D
 Acq On : 20 Jun 2022 20:45
 Operator : CG\JU
 Sample : N3373-06DL 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-6NDL

Quant Time: Jun 21 01:20:30 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	130867	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	423525	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	247251	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	456001	20.000	ng	0.00
76) Chrysene-d12	13.939	240	267022	20.000	ng	#-0.01
86) Perylene-d12	15.392	264	215877	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	62984	7.666	ng	0.00
7) Phenol-d6	6.428	99	86069	8.582	ng	-0.01
23) Nitrobenzene-d5	7.333	82	51942	5.710	ng	-0.02
42) 2,4,6-Tribromophenol	10.604	330	5959	2.040	ng	0.00
45) 2-Fluorobiphenyl	9.127	172	108125	6.126	ng	-0.01
79) Terphenyl-d14	12.880	244	118469	7.709	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.327	178	688743	30.113	ng	98
73) Carbazole	11.533	167	76025	3.616	ng	99
75) Fluoranthene	12.521	202	1190276	49.785	ng	97
78) Pyrene	12.745	202	701581	35.315	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062022\
Data File : BF128888.D
Acq On : 20 Jun 2022 20:45
Operator : CG\JU
Sample : N3373-06DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SC-6NDL

Quant Time: Jun 21 01:20:30 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

