

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131131.D
 Acq On : 10 Nov 2022 19:37
 Operator : CG\JU
 Sample : N5491-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

Quant Time: Nov 11 00:50:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	73334	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	277781	20.000	ng	#-0.01
39) Acenaphthene-d10	9.933	164	144677	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	242113	20.000	ng	0.00
76) Chrysene-d12	14.074	240	151704	20.000	ng	0.00
86) Perylene-d12	15.568	264	162012	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	499596	115.856	ng	0.00
7) Phenol-d6	6.516	99	634001	112.458	ng	0.00
23) Nitrobenzene-d5	7.451	82	428110	69.536	ng	-0.01
42) 2,4,6-Tribromophenol	10.727	330	147470	98.629	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	767856	74.090	ng	0.00
79) Terphenyl-d14	13.016	244	558071	64.499	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.080	276	2938	4.069	ng	# 92
92) Benzo(g,h,i)perylene	17.533	276	3151	4.617	ng	# 94

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

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