

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131144.D
 Acq On : 11 Nov 2022 02:17
 Operator : CG\JU
 Sample : N5495-04 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-31A

Quant Time: Nov 11 03:13:43 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	6.892	152	57831	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	180102	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	94470	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	202505	20.000	ng	0.00
76) Chrysene-d12	14.080	240	150880	20.000	ng	0.00
86) Perylene-d12	15.580	264	108357	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	157597	46.344	ng	0.00
7) Phenol-d6	6.516	99	184623	41.527	ng	0.00
23) Nitrobenzene-d5	7.451	82	114418	28.664	ng	-0.01
42) 2,4,6-Tribromophenol	10.727	330	41106	42.103	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	186172	27.510	ng	0.00
79) Terphenyl-d14	13.016	244	213130	24.767	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.092	276	7954	4.690	ng	Qvalue # 93
92) Benzo(g,h,i)perylene	17.551	276	8926	5.431	ng	# 85

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

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