

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131123.D
 Acq On : 10 Nov 2022 14:59
 Operator : CG\JU
 Sample : N5494-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-19

Quant Time: Nov 11 00:48:37 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	6.886	152	70692	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	270402	20.000	ng	-0.01
39) Acenaphthene-d10	9.933	164	144867	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	254212	20.000	ng	0.00
76) Chrysene-d12	14.074	240	156978	20.000	ng	0.00
86) Perylene-d12	15.574	264	162496	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	493108	118.625	ng	0.00
7) Phenol-d6	6.516	99	580659	106.846	ng	0.00
23) Nitrobenzene-d5	7.451	82	384323	64.128	ng	-0.01
42) 2,4,6-Tribromophenol	10.727	330	145258	97.022	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	676830	65.221	ng	0.00
79) Terphenyl-d14	13.016	244	581972	65.002	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.086	276	4099	4.148	ng	Qvalue 95
92) Benzo(g,h,i)perylene	17.545	276	4397	4.716	ng	# 86

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

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