

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110522\
 Data File : BF131031.D
 Acq On : 05 Nov 2022 21:32
 Operator : CG\JU
 Sample : N5378-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-16-20221028-10.5-11.5

Quant Time: Nov 07 00:52:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 00:46:02 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	84296	20.000	ng	0.00
21) Naphthalene-d8	8.187	136	309695	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	166846	20.000	ng	0.00
64) Phenanthrene-d10	11.439	188	289295	20.000	ng	0.00
76) Chrysene-d12	14.086	240	167513	20.000	ng	0.00
86) Perylene-d12	15.586	264	140276	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	609156	125.278	ng	0.00
7) Phenol-d6	6.528	99	773413	122.400	ng	0.00
23) Nitrobenzene-d5	7.469	82	519794	101.206	ng	0.00
42) 2,4,6-Tribromophenol	10.739	330	195360	139.095	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	809648	73.367	ng	0.00
79) Terphenyl-d14	13.027	244	733985	80.292	ng	0.00
Target Compounds						
						Qvalue
78) Pyrene	12.886	202	33370	2.361	ng	97
88) Benzo(b)fluoranthene	15.139	252	20350	2.244	ng	95
90) Benzo(a)pyrene	15.516	252	17744	2.434	ng	98

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

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