

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110722\
 Data File : BF131052.D
 Acq On : 07 Nov 2022 20:03
 Operator : CG\JU
 Sample : N5415-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-20

Quant Time: Nov 08 02:18:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 01:40:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	89711	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	359020	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	161305	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	246506	20.000	ng	0.00
76) Chrysene-d12	14.080	240	158356	20.000	ng	#-0.01
86) Perylene-d12	15.586	264	162196	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	498162	89.541	ng	0.00
7) Phenol-d6	6.522	99	612152	90.147	ng	0.00
23) Nitrobenzene-d5	7.463	82	398731	51.480	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	137793	77.589	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	664197	56.116	ng	0.00
79) Terphenyl-d14	13.021	244	437843	44.329	ng	-0.01
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.092	276	4438	3.625	ng	# 93
92) Benzo(g,h,i)perylene	17.551	276	4276	3.741	ng	# 72

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

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