

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110722\
 Data File : BF131051.D
 Acq On : 07 Nov 2022 19:29
 Operator : CG\JU
 Sample : N5415-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-17

Quant Time: Nov 08 02:17:48 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	91215	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	364995	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	195907	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	310115	20.000	ng	0.00
76) Chrysene-d12	14.080	240	219687	20.000	ng	-0.01
86) Perylene-d12	15.586	264	223128	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	514674	90.984	ng	0.00
7) Phenol-d6	6.522	99	617834	89.484	ng	0.00
23) Nitrobenzene-d5	7.463	82	384058	48.774	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	144737	67.105	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	742131	51.626	ng	0.00
79) Terphenyl-d14	13.027	244	601549	43.901	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.098	276	8473	3.743	ng	Qvalue # 92
92) Benzo(g,h,i)perylene	17.556	276	7353	3.829	ng	# 86

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

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