

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110722\
 Data File : BF131047.D
 Acq On : 07 Nov 2022 17:06
 Operator : CG\JU
 Sample : N5415-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-14

Quant Time: Nov 08 02:16:54 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	77378	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	365975	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	184569	20.000	ng	0.00
64) Phenanthrene-d10	11.439	188	259796	20.000	ng	# 0.00
76) Chrysene-d12	14.080	240	156512	20.000	ng	#-0.01
86) Perylene-d12	15.580	264	153224	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	459783	95.815	ng	0.00
7) Phenol-d6	6.528	99	532410	90.901	ng	0.00
23) Nitrobenzene-d5	7.463	82	365876	46.340	ng	0.00
42) 2,4,6-Tribromophenol	10.739	330	138494	68.155	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	735125	54.280	ng	0.00
79) Terphenyl-d14	13.027	244	648777	66.459	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.092	276	2324	3.490	ng	# 92
92) Benzo(g,h,i)perylene	17.551	276	2152	3.578	ng	# 74

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

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