

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110222\
 Data File : BF130936.D
 Acq On : 02 Nov 2022 18:16
 Operator : CG\JU
 Sample : N5395-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-12-103122

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/03/2022
 Supervised By : Jagrut Upadhyay 11/03/2022

Quant Time: Nov 03 03:44:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 03 03:37:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.922	152	75398	20.000	ng	0.00
21) Naphthalene-d8	8.204	136	292978	20.000	ng	-0.01
39) Acenaphthene-d10	9.969	164	152432	20.000	ng	0.00
64) Phenanthrene-d10	11.463	188	254246	20.000	ng	# 0.00
76) Chrysene-d12	14.110	240	136299	20.000	ng	-0.01
86) Perylene-d12	15.621	264	139612	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.551	112	615756	141.580	ng	0.01
7) Phenol-d6	6.545	99	781717	138.314	ng	0.00
23) Nitrobenzene-d5	7.487	82	526340	108.328	ng	0.00
42) 2,4,6-Tribromophenol	10.763	330	189508	147.687	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	835750	82.893	ng	0.00
79) Terphenyl-d14	13.051	244	614551	82.622	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	12.674	202	33865	2.429	ng	97
78) Pyrene	12.904	202	32978	2.868	ng	98
88) Benzo(b)fluoranthene	15.174	252	19262m	2.134	ng	
90) Benzo(a)pyrene	15.551	252	15462	2.131	ng	# 89

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

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