

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110222\
 Data File : BF130939.D
 Acq On : 02 Nov 2022 20:01
 Operator : CG\JU
 Sample : N5395-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-3-103122

Manual Integrations
 APPROVED

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

Quant Time: Nov 03 03:46:19 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	72012	20.000	ng	0.00
21) Naphthalene-d8	8.204	136	275412	20.000	ng	-0.01
39) Acenaphthene-d10	9.969	164	142566	20.000	ng	0.00
64) Phenanthrene-d10	11.457	188	233161	20.000	ng	-0.01
76) Chrysene-d12	14.110	240	144916	20.000	ng	-0.01
86) Perylene-d12	15.621	264	146758	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.545	112	483481	116.393	ng	0.00
7) Phenol-d6	6.545	99	607019	112.454	ng	0.00
23) Nitrobenzene-d5	7.487	82	415687	91.011	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	150329	125.262	ng	-0.01
45) 2-Fluorobiphenyl	9.286	172	665863	70.613	ng	0.00
79) Terphenyl-d14	13.051	244	525329	66.428	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	7.316	107	19507	3.793	ng	84
71) Phenanthrene	11.480	178	37565	3.122	ng	99
75) Fluoranthene	12.674	202	76208	5.960	ng	98
78) Pyrene	12.904	202	63082	5.160	ng	99
81) Benzo(a)anthracene	14.098	228	32366	3.369	ng	96
83) Chrysene	14.133	228	30107	3.249	ng	97
88) Benzo(b)fluoranthene	15.168	252	42751m	4.506	ng	
90) Benzo(a)pyrene	15.551	252	28937	3.794	ng	95
92) Benzo(g,h,i)perylene	17.615	276	18918	2.334	ng	# 90

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

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