

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081722\
 Data File : BF129843.D
 Acq On : 17 Aug 2022 13:58
 Operator : CG\JU
 Sample : N4189-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB17

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/18/2022
 Supervised By : Jagrut Upadhyay 08/18/2022

Quant Time: Aug 17 16:47:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 15 02:28:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	139583	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	536295	20.000	ng	0.00
39) Acenaphthene-d10	9.863	164	248664	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	309191	20.000	ng	0.00
76) Chrysene-d12	14.004	240	269688	20.000	ng	0.00
86) Perylene-d12	15.474	264	228742	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	742665	88.093	ng	0.00
7) Phenol-d6	6.481	99	922056	87.338	ng	0.00
23) Nitrobenzene-d5	7.392	82	580761	57.970	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	156350	62.647	ng	0.00
45) 2-Fluorobiphenyl	9.181	172	1049144	64.080	ng	0.00
79) Terphenyl-d14	12.939	244	750506	46.305	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	12.569	202	97598	5.523	ng	98
78) Pyrene	12.798	202	92011	3.926	ng	99
81) Benzo(a)anthracene	13.992	228	50592	2.727	ng	96
83) Chrysene	14.027	228	49638	2.858	ng	96
88) Benzo(b)fluoranthene	15.045	252	66883m	4.270	ng	
90) Benzo(a)pyrene	15.415	252	39020	3.163	ng	93

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

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