

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011022\
 Data File : BF126563.D
 Acq On : 10 Jan 2022 13:55
 Operator : JU/CG
 Sample : N1047-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FM-E7-01-2.5

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/11/2022
 Supervised By : Jagrut Upadhyay 01/11/2022

Quant Time: Jan 10 17:16:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 10 14:50:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	93780	20.000	ng	0.00
21) Naphthalene-d8	8.051	136	393294	20.000	ng	0.00
39) Acenaphthene-d10	9.810	164	186675	20.000	ng	0.00
64) Phenanthrene-d10	11.292	188	286801	20.000	ng	0.00
76) Chrysene-d12	13.939	240	169744	20.000	ng	0.00
86) Perylene-d12	15.392	264	178701	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	620622	95.036	ng	0.00
7) Phenol-d6	6.404	99	797528	94.291	ng	0.00
23) Nitrobenzene-d5	7.334	82	499745	65.244	ng	0.00
42) 2,4,6-Tribromophenol	10.598	330	164528	114.056	ng	0.00
45) 2-Fluorobiphenyl	9.133	172	742082	69.954	ng	0.00
79) Terphenyl-d14	12.886	244	575816	65.909	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	12.510	202	63676	4.758	ng	97
78) Pyrene	12.733	202	62392	4.349	ng	98
81) Benzo(a)anthracene	13.927	228	29898	3.063	ng	95
83) Chrysene	13.963	228	26207	2.617	ng	95
88) Benzo(b)fluoranthene	14.974	252	38795m	3.644	ng	
90) Benzo(a)pyrene	15.333	252	30393	3.405	ng	98

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

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