

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042922\
 Data File : BF127998.D
 Acq On : 29 Apr 2022 16:48
 Operator : CG\JU
 Sample : N2616-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-18D

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/02/2022
 Supervised By : Jagrut Upadhyay 05/02/2022

Quant Time: Apr 29 17:58:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 28 05:13:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.922	152	105373	20.000	ng	-0.01
21) Naphthalene-d8	8.204	136	408622	20.000	ng	-0.01
39) Acenaphthene-d10	9.963	164	236662	20.000	ng	-0.01
64) Phenanthrene-d10	11.457	188	421153	20.000	ng	-0.01
76) Chrysene-d12	14.098	240	238578	20.000	ng	#-0.02
86) Perylene-d12	15.610	264	220101	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.546	112	571957	86.455	ng	0.00
7) Phenol-d6	6.546	99	759413	88.466	ng	-0.02
23) Nitrobenzene-d5	7.487	82	570594	65.664	ng	-0.02
42) 2,4,6-Tribromophenol	10.757	330	224971	94.430	ng	-0.01
45) 2-Fluorobiphenyl	9.281	172	938586	67.093	ng	-0.01
79) Terphenyl-d14	13.045	244	886492	67.944	ng	-0.01
Target Compounds						Qvalue
78) Pyrene	12.898	202	46001	2.389	ng	99
88) Benzo(b)fluoranthene	15.157	252	33248m	2.405	ng	
90) Benzo(a)pyrene	15.539	252	24152	2.106	ng	97

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

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