

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
 Data File : BF128915.D
 Acq On : 21 Jun 2022 23:51
 Operator : CG\JU
 Sample : N3420-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MMC-CON-0620

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/22/2022
 Supervised By : Jagrut Upadhyay 06/22/2022

Quant Time: Jun 22 01:57:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 01:02:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.751	152	95918	20.000	ng	0.00
21) Naphthalene-d8	8.034	136	343755	20.000	ng	0.00
39) Acenaphthene-d10	9.792	164	171327	20.000	ng	0.00
64) Phenanthrene-d10	11.275	188	258779	20.000	ng	0.00
76) Chrysene-d12	13.921	240	205866	20.000	ng	# 0.00
86) Perylene-d12	15.362	264	148880	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	77512	12.872	ng	0.00
7) Phenol-d6	6.410	99	363681	49.476	ng	0.00
23) Nitrobenzene-d5	7.322	82	384880	52.129	ng	0.00
42) 2,4,6-Tribromophenol	10.586	330	997	0.493	ng	0.00
45) 2-Fluorobiphenyl	9.116	172	706717	57.782	ng	0.00
79) Terphenyl-d14	12.863	244	614907	51.902	ng	0.00
Target Compounds						
52) Acenaphthene	9.822	154	21626	2.211	ng	Qvalue 97
71) Phenanthrene	11.298	178	57709	4.446	ng	99
75) Fluoranthene	12.492	202	51827	3.820	ng	96
78) Pyrene	12.721	202	46111	3.011	ng	97
83) Chrysene	13.945	228	25032	2.046	ng	# 95
88) Benzo(b)fluoranthene	14.945	252	22466m	2.362	ng	

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

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