

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061522\
 Data File : BF128826.D
 Acq On : 15 Jun 2022 15:04
 Operator : CG\JU
 Sample : N3337-01DL2 20X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-02-061322DL2

Manual Integrations
 APPROVED

Reviewed By :Jagrut
 Upadhyay

06/16/2022
 Supervised By :mohammad
 ahmed

Quant Time: Jun 15 15:35:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 14 14:53:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	86581	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	314659	20.000	ng	# 0.00
39) Acenaphthene-d10	9.822	164	158161	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	260502	20.000	ng	0.00
76) Chrysene-d12	13.957	240	177021	20.000	ng	# 0.00
86) Perylene-d12	15.415	264	135558	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	24320	4.474	ng	0.00
7) Phenol-d6	6.439	99	27446	4.136	ng	0.00
23) Nitrobenzene-d5	7.339	82	17750	2.626	ng	-0.01
42) 2,4,6-Tribromophenol	10.616	330	4830	2.585	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	36896	3.268	ng	0.00
79) Terphenyl-d14	12.898	244	31771	3.119	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.086	128	57573	3.565	ng	99
52) Acenaphthene	9.857	154	47331	5.243	ng	97
55) Dibenzofuran	10.027	168	54689	4.215	ng	98
58) Fluorene	10.369	166	65845	6.544	ng	96
71) Phenanthrene	11.339	178	487639	37.321	ng	98
72) Anthracene	11.386	178	160557	12.415	ng	97
73) Carbazole	11.545	167	25076	2.088	ng	98
75) Fluoranthene	12.533	202	638324	46.736	ng	95
78) Pyrene	12.763	202	529998	40.242	ng	99
81) Benzo(a)anthracene	13.951	228	235021	21.299	ng	99
83) Chrysene	13.986	228	187908	17.862	ng	97
87) Indeno(1,2,3-cd)pyrene	16.856	276	75318	9.218	ng	# 87
88) Benzo(b)fluoranthene	14.992	252	204378m	23.595	ng	
89) Benzo(k)fluoranthene	15.021	252	74492m	9.142	ng	
90) Benzo(a)pyrene	15.357	252	153102m	22.489	ng	
91) Dibenzo(a,h)anthracene	16.868	278	20340m	3.001	ng	
92) Benzo(g,h,i)perylene	17.298	276	68154	10.147	ng	# 89

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061522\
Data File : BF128826.D
Acq On : 15 Jun 2022 15:04
Operator : CG\JU
Sample : N3337-01DL2 20X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-02-061322DL2

Quant Time: Jun 15 15:35:52 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 14 14:53:53 2022
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Jagrut
Upadhyay

06/16/2022
Supervised By :mohammad
ahmed

06/22/2022

