

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080322\
 Data File : BF129554.D
 Acq On : 03 Aug 2022 22:44
 Operator : CG\JU
 Sample : N3930-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-99

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/04/2022
 Supervised By : Jagrut Upadhyay 08/05/2022

Quant Time: Aug 04 02:21:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	128600	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	447515	20.000	ng	-0.01
39) Acenaphthene-d10	9.904	164	184837	20.000	ng	-0.01
64) Phenanthrene-d10	11.398	188	292210	20.000	ng	0.00
76) Chrysene-d12	14.039	240	319864	20.000	ng	-0.01
86) Perylene-d12	15.527	264	274253	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	460124	58.285	ng	0.00
7) Phenol-d6	6.510	99	541271	54.548	ng	-0.01
23) Nitrobenzene-d5	7.428	82	357762	43.640	ng	-0.02
42) 2,4,6-Tribromophenol	10.698	330	97917	55.100	ng	-0.01
45) 2-Fluorobiphenyl	9.222	172	569243	47.641	ng	-0.01
79) Terphenyl-d14	12.980	244	591276	34.874	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	12.610	202	63079	3.903	ng	98
78) Pyrene	12.839	202	76858	2.873	ng	97
81) Benzo(a)anthracene	14.027	228	43725	2.001	ng	94
83) Chrysene	14.068	228	44238	2.148	ng	98
88) Benzo(b)fluoranthene	15.086	252	54895m	3.018	ng	
90) Benzo(a)pyrene	15.463	252	35297	2.390	ng	# 94

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

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