

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072822\
 Data File : BF129425.D
 Acq On : 29 Jul 2022 00:09
 Operator : CG\JU
 Sample : N3893-20
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-19D

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 07/29/2022
 Supervised By :Jagrut Upadhyay 07/29/2022

Quant Time: Jul 29 03:26:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 17:24:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	193004	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	745895	20.000	ng	# 0.00
39) Acenaphthene-d10	9.922	164	358834	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	487110	20.000	ng	0.00
76) Chrysene-d12	14.063	240	369239	20.000	ng	0.00
86) Perylene-d12	15.557	264	336447	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1006876	84.983	ng	0.01
7) Phenol-d6	6.522	99	1302095	87.434	ng	0.00
23) Nitrobenzene-d5	7.445	82	820921	60.079	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	270720	78.471	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	1480362	63.819	ng	0.00
79) Terphenyl-d14	12.998	244	1121659	57.310	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	12.628	202	106705	3.961	ng	98
78) Pyrene	12.857	202	153294	4.965	ng	98
83) Chrysene	14.086	228	86100	3.622	ng	98
88) Benzo(b)fluoranthene	15.110	252	95550m	4.281	ng	
90) Benzo(a)pyrene	15.486	252	59604	3.290	ng	# 93

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

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