

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
 Data File : BF129441.D
 Acq On : 29 Jul 2022 14:34
 Operator : CG\JU
 Sample : N3896-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-73

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/01/2022
 Supervised By : Jagrut Upadhyay 08/01/2022

Quant Time: Jul 29 18:21:29 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	184427	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	736148	20.000	ng	0.00
39) Acenaphthene-d10	9.927	164	382077	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	628060	20.000	ng	# 0.00
76) Chrysene-d12	14.062	240	364993	20.000	ng	0.00
86) Perylene-d12	15.557	264	385156	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1109290	97.981	ng	0.02
7) Phenol-d6	6.522	99	1434501	100.805	ng	0.00
23) Nitrobenzene-d5	7.451	82	912427	67.660	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	378146	102.942	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1629809	65.988	ng	0.00
79) Terphenyl-d14	13.004	244	1441070	74.486	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	12.633	202	87111	2.508	ng	99
78) Pyrene	12.863	202	93958	3.078	ng	99
88) Benzo(b)fluoranthene	15.115	252	59659m	2.335	ng	

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

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