

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
 Data File : BF129438.D
 Acq On : 29 Jul 2022 12:57
 Operator : CG\JU
 Sample : N3896-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-74

Manual Integrations
 APPROVED

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

Quant Time: Jul 29 18:20:26 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	6.887	152	182984	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	733134	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	383254	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	604149	20.000	ng	# 0.00
76) Chrysene-d12	14.063	240	371245	20.000	ng	# 0.00
86) Perylene-d12	15.562	264	376324	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1116537	99.399	ng	0.02
7) Phenol-d6	6.522	99	1425866	100.988	ng	0.00
23) Nitrobenzene-d5	7.451	82	925963	68.946	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	374704	101.692	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1663369	67.139	ng	0.00
79) Terphenyl-d14	13.004	244	1454291	73.904	ng	0.00
Target Compounds						
75) Fluoranthene	12.633	202	99181	2.968	ng	99
78) Pyrene	12.863	202	95688	3.082	ng	97

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

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