

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072822\
 Data File : BF129424.D
 Acq On : 28 Jul 2022 23:36
 Operator : CG\JU
 Sample : N3893-19
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-19

Manual Integrations
 APPROVED

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

Quant Time: Jul 29 02:10:22 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	185541	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	723679	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	348692	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	480179	20.000	ng	0.00
76) Chrysene-d12	14.063	240	354162	20.000	ng	0.00
86) Perylene-d12	15.551	264	324842	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1085492	95.303	ng	0.01
7) Phenol-d6	6.522	99	1375852	96.103	ng	0.00
23) Nitrobenzene-d5	7.445	82	875004	66.003	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	289184	86.261	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	1550882	68.804	ng	0.00
79) Terphenyl-d14	12.998	244	1177269	62.712	ng	0.00
Target Compounds						
75) Fluoranthene	12.627	202	123970	4.668	ng	99
78) Pyrene	12.863	202	181781	6.138	ng	97
83) Chrysene	14.086	228	101635	4.458	ng	97
88) Benzo(b)fluoranthene	15.110	252	101841m	4.726	ng	
90) Benzo(a)pyrene	15.486	252	71111	4.065	ng	94

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

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