

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080422\
 Data File : BF129591.D
 Acq On : 05 Aug 2022 03:04
 Operator : CG\JU
 Sample : N3930-15
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-107

Manual Integrations
 APPROVED

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

Quant Time: Aug 05 03:31:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 02:05:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	142072	20.000	ng	0.00
21) Naphthalene-d8	8.140	136	498913	20.000	ng	0.00
39) Acenaphthene-d10	9.898	164	211719	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	313040	20.000	ng	0.00
76) Chrysene-d12	14.033	240	260418	20.000	ng	0.00
86) Perylene-d12	15.516	264	210345	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	739584	84.801	ng	0.00
7) Phenol-d6	6.510	99	882645	80.516	ng	0.00
23) Nitrobenzene-d5	7.422	82	532986	58.317	ng	-0.01
42) 2,4,6-Tribromophenol	10.686	330	156234	76.754	ng	-0.01
45) 2-Fluorobiphenyl	9.216	172	899648	65.734	ng	0.00
79) Terphenyl-d14	12.969	244	791485	57.339	ng	-0.01
Target Compounds						
75) Fluoranthene	12.604	202	57194	3.303	ng	Qvalue 98
78) Pyrene	12.833	202	60946	2.799	ng	96
81) Benzo(a)anthracene	14.022	228	36882	2.073	ng	93
83) Chrysene	14.057	228	55747	3.325	ng	98
88) Benzo(b)fluoranthene	15.080	252	89553m	6.418	ng	
90) Benzo(a)pyrene	15.451	252	30389	2.683	ng	# 92

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

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