

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110722\
 Data File : BF131054.D
 Acq On : 07 Nov 2022 21:10
 Operator : CG\JU
 Sample : N5415-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-21

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/08/2022
 Supervised By : Jagrut Upadhyay 11/08/2022

Quant Time: Nov 08 02:18:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 01:40:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	76313	20.000	ng	0.00
21) Naphthalene-d8	8.187	136	319222	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	137436	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	173751	20.000	ng	0.00
76) Chrysene-d12	14.086	240	193230	20.000	ng	0.00
86) Perylene-d12	15.586	264	160443	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	616540	130.275	ng	0.00
7) Phenol-d6	6.528	99	776032	134.345	ng	0.00
23) Nitrobenzene-d5	7.469	82	533276	77.435	ng	0.00
42) 2,4,6-Tribromophenol	10.739	330	127823	84.476	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	766502	76.007	ng	0.00
79) Terphenyl-d14	13.027	244	616073	51.116	ng	0.00
Target Compounds						
75) Fluoranthene	12.651	202	22061	2.124	ng	Qvalue 97
87) Indeno(1,2,3-cd)pyrene	17.098	276	5951	3.733	ng	# 85
88) Benzo(b)fluoranthene	15.139	252	23678m	2.031	ng	
92) Benzo(g,h,i)perylene	17.562	276	5934	3.882	ng	# 86

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

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