

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032321\
 Data File : BF123426.D
 Acq On : 24 Mar 2021 7:00
 Operator : JU/CG
 Sample : M1712-03
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EXTERNAL-PAD-2

Manual Integrations
 APPROVED

mohammad
 3/24/2021 1:33:35 PM

Quant Time: Mar 24 07:38:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 23 15:58:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	281410	20.00	ng	0.00
21) Naphthalene-d8	8.198	136	1075562	20.00	ng	# 0.00
39) Acenaphthene-d10	9.957	164	588363	20.00	ng	0.00
64) Phenanthrene-d10	11.451	188	897322	20.00	ng	0.00
76) Chrysene-d12	14.098	240	437316	20.00	ng	0.00
86) Perylene-d12	15.604	264	462719	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	1258458	73.16	ng	0.02
7) Phenol-d6	6.534	99	1638819	71.57	ng	0.00
23) Nitrobenzene-d5	7.475	82	1133353	56.87	ng	0.00
42) 2,4,6-Tribromophenol	10.745	330	368867	56.08	ng	-0.01
45) 2-Fluorobiphenyl	9.275	172	2094684	50.86	ng	0.00
79) Terphenyl-d14	13.045	244	1597910	61.93	ng	0.00
Target Compounds						
78) Pyrene	12.892	202	126993	3.37	ng	Qvalue 97
83) Chrysene	14.121	228	76997	2.61	ng	97
87) Indeno(1,2,3-cd)pyrene	17.115	276	116249	4.04	ng	98
88) Benzo(b)fluoranthene	15.157	252	124522m	3.90	ng	
92) Benzo(g,h,i)perylene	17.574	276	184073	7.53	ng	99

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

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