

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
 Data File : BF129731.D
 Acq On : 11 Aug 2022 23:51
 Operator : CG\JU
 Sample : N4130-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB03

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/12/2022
 Supervised By : Jagrut Upadhyay 08/12/2022

Quant Time: Aug 12 01:26:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 11 23:50:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	148810	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	556761	20.000	ng	# 0.00
39) Acenaphthene-d10	9.869	164	247571	20.000	ng	-0.01
64) Phenanthrene-d10	11.357	188	351826	20.000	ng	#-0.01
76) Chrysene-d12	14.010	240	278794	20.000	ng	-0.01
86) Perylene-d12	15.480	264	233650	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	785589	85.997	ng	0.00
7) Phenol-d6	6.487	99	966637	84.186	ng	0.00
23) Nitrobenzene-d5	7.399	82	601984	59.023	ng	-0.01
42) 2,4,6-Tribromophenol	10.663	330	176662	74.221	ng	-0.01
45) 2-Fluorobiphenyl	9.187	172	1010761	63.157	ng	-0.01
79) Terphenyl-d14	12.945	244	769092	52.044	ng	-0.01
Target Compounds						Qvalue
75) Fluoranthene	12.575	202	69091	3.551	ng	98
78) Pyrene	12.804	202	61839	2.652	ng	98
84) Bis(2-ethylhexyl)phtha...	13.969	149	82017	6.161	ng	99
88) Benzo(b)fluoranthene	15.051	252	42037m	2.712	ng	
90) Benzo(a)pyrene	15.416	252	25239	2.006	ng	# 90

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

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