

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082823\
 Data File : BF134975.D
 Acq On : 28 Aug 2023 18:41
 Operator : CG\JU
 Sample : 04110-10DL 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-11(1-2)DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/29/2023
 Supervised By :mohammad ahmed 08/29/2023

Quant Time: Aug 29 03:51:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 28 02:45:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.793	152	106085	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	383960	20.000	ng	0.00
39) Acenaphthene-d10	9.828	164	172487	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	287983	20.000	ng	-0.01
76) Chrysene-d12	13.980	240	257775	20.000	ng	0.00
86) Perylene-d12	15.451	264	224479	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	116276	17.478	ng	0.00
7) Phenol-d6	6.422	99	263379	30.983	ng	-0.02
23) Nitrobenzene-d5	7.345	82	199703	28.660	ng	-0.02
42) 2,4,6-Tribromophenol	10.616	330	12247	6.029	ng	-0.01
45) 2-Fluorobiphenyl	9.145	172	343737	30.807	ng	-0.01
79) Terphenyl-d14	12.916	244	329789	17.596	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.092	128	89153	4.650	ng	99
52) Acenaphthene	9.857	154	47126	4.777	ng	98
55) Dibenzofuran	10.034	168	31241	2.175	ng	# 93
58) Fluorene	10.375	166	28203	2.589	ng	99
71) Phenanthrene	11.345	178	373462	24.748	ng	99
72) Anthracene	11.392	178	89315	5.776	ng	98
73) Carbazole	11.551	167	53061	4.165	ng	97
75) Fluoranthene	12.539	202	608039	42.524	ng	99
78) Pyrene	12.769	202	550485	20.141	ng	100
81) Benzo(a)anthracene	13.969	228	371183	21.105	ng	97
83) Chrysene	14.004	228	366730	21.134	ng	97
87) Indeno(1,2,3-cd)pyrene	16.939	276	159393	9.934	ng	96
88) Benzo(b)fluoranthene	15.021	252	445916m	35.209	ng	
89) Benzo(k)fluoranthene	15.051	252	160384m	12.896	ng	
90) Benzo(a)pyrene	15.392	252	311024	24.600	ng	98
91) Dibenzo(a,h)anthracene	16.951	278	41406	3.189	ng	# 90
92) Benzo(g,h,i)perylene	17.398	276	151093	11.369	ng	97

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

