

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072723\
 Data File : BF134503.D
 Acq On : 27 Jul 2023 13:38
 Operator : CG\JU
 Sample : 03640-31DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP11DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/28/2023
 Supervised By :mohammad ahmed 07/28/2023

Quant Time: Jul 27 13:58:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 27 12:16:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	53265	20.000	ng	0.00
21) Naphthalene-d8	8.098	136	214099	20.000	ng	# 0.00
39) Acenaphthene-d10	9.857	164	111368	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	218390	20.000	ng	0.00
76) Chrysene-d12	14.016	240	138082	20.000	ng	0.00
86) Perylene-d12	15.510	264	158223	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	90136	27.959	ng	0.00
7) Phenol-d6	6.469	99	59980	15.185	ng	0.00
23) Nitrobenzene-d5	7.387	82	40455	9.560	ng	0.00
42) 2,4,6-Tribromophenol	10.651	330	23238	14.686	ng	-0.01
45) 2-Fluorobiphenyl	9.175	172	89049	10.497	ng	0.00
79) Terphenyl-d14	12.945	244	83759	10.589	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.122	128	42138	3.955	ng	98
37) 2-Methylnaphthalene	8.816	142	20281	2.739	ng	95
38) 1-Methylnaphthalene	8.916	142	17975	2.617	ng	95
55) Dibenzofuran	10.063	168	24899	2.396	ng	# 96
58) Fluorene	10.410	166	37769	4.504	ng	99
71) Phenanthrene	11.381	178	468882	41.092	ng	100
72) Anthracene	11.428	178	106471	9.041	ng	99
73) Carbazole	11.592	167	25131	2.266	ng	99
75) Fluoranthene	12.575	202	547921	41.202	ng	99
78) Pyrene	12.810	202	460978	41.986	ng	99
81) Benzo(a)anthracene	14.004	228	235480	24.443	ng	99
83) Chrysene	14.039	228	186167	20.204	ng	99
87) Indeno(1,2,3-cd)pyrene	17.045	276	93113	9.008	ng	99
88) Benzo(b)fluoranthene	15.069	252	240485m	24.295	ng	
89) Benzo(k)fluoranthene	15.098	252	87245m	8.962	ng	
90) Benzo(a)pyrene	15.445	252	178824	19.545	ng	# 97
91) Dibenzo(a,h)anthracene	17.045	278	24502	2.935	ng	93
92) Benzo(g,h,i)perylene	17.510	276	80676	9.412	ng	98

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

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