

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060722\
 Data File : BF128723.D
 Acq On : 07 Jun 2022 18:46
 Operator : CG\JU
 Sample : N3201-03DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 21500DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/08/2022
 Supervised By : Jagrut Upadhyay 06/08/2022

Quant Time: Jun 08 00:27:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 02 17:07:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	77803	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	275021	20.000	ng	# 0.00
39) Acenaphthene-d10	9.863	164	140899	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	211941	20.000	ng	# 0.00
76) Chrysene-d12	13.998	240	174901	20.000	ng	# 0.00
86) Perylene-d12	15.463	264	190525	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	35617	7.292	ng	0.00
7) Phenol-d6	6.498	99	42506	7.129	ng	0.00
23) Nitrobenzene-d5	7.381	82	30102	5.096	ng	-0.01
42) 2,4,6-Tribromophenol	10.657	330	8221	4.938	ng	0.00
45) 2-Fluorobiphenyl	9.181	172	42921	4.267	ng	-0.01
79) Terphenyl-d14	12.933	244	44457	4.417	ng	-0.01
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.822	142	38657	4.306	ng	100
38) 1-Methylnaphthalene	8.922	142	19922	2.297	ng	93
52) Acenaphthene	9.892	154	50202	6.242	ng	96
55) Dibenzofuran	10.069	168	65006	5.624	ng	95
58) Fluorene	10.410	166	86230	9.620	ng	99
71) Phenanthrene	11.375	178	313488	29.489	ng	99
72) Anthracene	11.428	178	40067	3.808	ng	98
75) Fluoranthene	12.563	202	221898	19.969	ng	96
78) Pyrene	12.792	202	147541	11.338	ng	99
81) Benzo(a)anthracene	13.986	228	43057	3.949	ng	96
83) Chrysene	14.022	228	48142	4.632	ng	99
88) Benzo(b)fluoranthene	15.033	252	31791m	2.611	ng	

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

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