

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101923\
 Data File : BF135895.D
 Acq On : 19 Oct 2023 19:49
 Operator : CG\JU
 Sample : 04704-04DL 25X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-3(10-15)DL

Quant Time: Oct 20 02:04:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 17 01:22:00 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/20/2023
 Supervised By :mohammad ahmed 10/20/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							10/20/2023
1) 1,4-Dichlorobenzene-d4	6.734	152	58049	20.000	ng	-0.01	Supervised By :mohammad ahmed
21) Naphthalene-d8	8.016	136	203534	20.000	ng	0.00	
39) Acenaphthene-d10	9.774	164	88532	20.000	ng	-0.01	
64) Phenanthrene-d10	11.269	188	147628	20.000	ng	-0.01	
76) Chrysene-d12	13.939	240	116014	20.000	ng	-0.01	
86) Perylene-d12	15.415	264	97445	20.000	ng	0.00	10/20/2023
System Monitoring Compounds							
5) 2-Fluorophenol	5.416	112	12120	3.357	ng	0.04	
7) Phenol-d6	6.434	99	16439	3.649	ng	0.04	
23) Nitrobenzene-d5	7.322	82	9081	2.454	ng	0.00	
42) 2,4,6-Tribromophenol	10.580	330	2727	3.210	ng	0.00	
45) 2-Fluorobiphenyl	9.092	172	19995	3.531	ng	-0.01	
79) Terphenyl-d14	12.863	244	18744	2.992	ng	-0.02	
Target Compounds						Qvalue	
31) Naphthalene	8.033	128	134577	13.556	ng	99	
37) 2-Methylnaphthalene	8.727	142	48922	7.480	ng	96	
38) 1-Methylnaphthalene	8.827	142	104626	17.343	ng	98	
46) 1,1'-Biphenyl	9.192	154	20700	3.052	ng	98	
49) Acenaphthylene	9.633	152	58414	7.222	ng	# 95	
52) Acenaphthene	9.804	154	15120	3.119	ng	96	
58) Fluorene	10.322	166	65023	11.354	ng	96	
71) Phenanthrene	11.292	178	294717	41.673	ng	100	
72) Anthracene	11.345	178	54461	7.432	ng	98	
75) Fluoranthene	12.492	202	178245	22.709	ng	97	
78) Pyrene	12.727	202	281350	30.518	ng	99	
81) Benzo(a)anthracene	13.927	228	105384	13.829	ng	97	
83) Chrysene	13.962	228	93815	12.711	ng	98	
87) Indeno(1,2,3-cd)pyrene	16.927	276	25579	4.818	ng	95	
88) Benzo(b)fluoranthene	14.986	252	60772m	11.143	ng		
89) Benzo(k)fluoranthene	15.004	252	19610m	3.704	ng		
90) Benzo(a)pyrene	15.356	252	61610	12.014	ng	98	
92) Benzo(g,h,i)perylene	17.386	276	27342	6.115	ng	# 89	

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

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