

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021020\
 Data File : BF118876.D
 Acq On : 10 Feb 2020 18:44
 Operator : CG/JU
 Sample : L1470-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PE-6-(SIDE-WALL)-NORTH

Manual Integrations
 APPROVED

mohammad
 2/13/2020 10:28:48 AM

Quant Time: Feb 11 00:50:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 03 16:26:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	245971	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	913418	20.00	ng	-0.01
39) Acenaphthene-d10	9.85	164	529124	20.00	ng	-0.01
64) Phenanthrene-d10	11.33	188	1052942	20.00	ng	0.00
76) Chrysene-d12	13.97	240	800487	20.00	ng	-0.01
87) Perylene-d12	15.42	264	865254	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1395219	107.37	ng	0.00
7) Phenol-d6	6.45	99	1715727	106.37	ng	0.00
23) Nitrobenzene-d5	7.37	82	1134717	74.05	ng	-0.01
42) 2,4,6-Tribromophenol	10.64	330	673475	106.38	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	2259751	76.35	ng	-0.01
79) Terphenyl-d14	12.92	244	2857203	73.32	ng	-0.01

Target Compounds				Qvalue		
31) Naphthalene	8.12	128	88290	2.113	ng	98
37) 2-Methylnaphthalene	8.80	142	66541	2.297	ng	97
38) 1-Methylnaphthalene	8.90	142	55003	2.018	ng	100
49) Acenaphthylene	9.71	152	149939	3.413	ng	98
50) Dimethylphthalate	9.56	163	78490	2.179	ng	99
52) Acenaphthene	9.88	154	60088	2.076	ng	99
58) Fluorene	10.39	166	119447	3.795	ng	98
71) Phenanthrene	11.36	178	1673353	32.300	ng	99
72) Anthracene	11.40	178	413169	8.017	ng	98
75) Fluoranthene	12.55	202	3103037	54.860	ng	98
78) Pyrene	12.77	202	2818418	51.323	ng	99
81) Benzo(a)anthracene	13.96	228	1576176	33.089	ng	97
83) Chrysene	14.00	228	1385945	28.990	ng	98
86) Indeno(1,2,3-cd)pyrene	16.86	276	964068	20.071	ng	96
88) Benzo(b)fluoranthene	15.00	252	1801221m	33.548	ng	
89) Benzo(k)fluoranthene	15.02	252	712820m	15.246	ng	
90) Benzo(a)pyrene	15.36	252	1387039	30.187	ng	99
91) Dibenzo(a,h)anthracene	16.86	278	255318	6.259	ng	94
92) Benzo(a,h,i)perylene	17.29	276	859213	21.903	ng	98

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

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