

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
 Data File : BP001655.D
 Acq On : 17 Jan 2020 13:58
 Operator : JU
 Sample : L1148-06DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-27-(4-6)DL

Manual Integrations
 APPROVED

mohammad
 1/21/2020 7:52:51 AM

Quant Time: Jan 18 03:07:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 21:52:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	53096	20.00	ng	0.00
21) Naphthalene-d8	10.41	136	211092	20.00	ng	0.00
39) Acenaphthene-d10	14.28	164	151225	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	317679	20.00	ng	0.00
76) Chrysene-d12	21.14	240	199991	20.00	ng	-0.01
87) Perylene-d12	23.38	264	187918	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	133291	49.18	ng	0.00
7) Phenol-d6	6.83	99	208977	48.24	ng	0.00
23) Nitrobenzene-d5	8.78	82	147559	30.55	ng	0.00
42) 2,4,6-Tribromophenol	15.78	330	80051	34.78	ng	-0.01
45) 2-Fluorobiphenyl	12.90	172	307312	29.88	ng	0.00
79) Terphenyl-d14	19.63	244	404204	36.97	ng	-0.01
Target Compounds						
49) Acenaphthylene	14.00	152	55388	3.988	ng	99
50) Dimethylphthalate	13.75	163	29559	2.373	ng	97
52) Acenaphthene	14.35	154	96972	11.273	ng	98
55) Dibenzofuran	14.68	168	141807	9.748	ng	98
58) Fluorene	15.34	166	178647	15.176	ng	99
71) Phenanthrene	17.09	178	960208	55.579	ng	99
72) Anthracene	17.17	178	155861	9.267	ng	99
73) Carbazole	17.45	167	64988	4.116	ng	98
75) Fluoranthene	19.07	202	635019	27.934	ng	99
78) Pyrene	19.42	202	509105	34.654	ng	98
81) Benzo(a)anthracene	21.13	228	172260m	12.368	ng	
83) Chrysene	21.18	228	161795	11.920	ng	95
86) Indeno(1,2,3-cd)pyrene	25.63	276	84605	5.325	ng	98
88) Benzo(b)fluoranthene	22.71	252	173796m	14.677	ng	
89) Benzo(k)fluoranthene	22.74	252	58914m	5.103	ng	
90) Benzo(a)pyrene	23.28	252	121260	10.785	ng	# 98
91) Dibenzo(a,h)anthracene	25.63	278	25177	2.170	ng	# 94
92) Benzo(g,h,i)perylene	26.31	276	88902	7.709	ng	98

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

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