

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090220\
 Data File : BP003186.D
 Acq On : 02 Sep 2020 14:46
 Operator : CG/JU
 Sample : L3848-23DL 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 LINDEN-BH-01-CDL

Manual Integrations
 APPROVED

mohammad
 9/3/2020 1:48:26 PM

Quant Time: Sep 08 08:38:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 14:37:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	205494	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	863124	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	542095	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1209539	20.00	ng	0.00
76) Chrysene-d12	21.15	240	1232280	20.00	ng	0.00
86) Perylene-d12	23.38	264	1312567	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	449800	38.70	ng	0.00
7) Phenol-d6	6.84	99	552664	37.93	ng	0.00
23) Nitrobenzene-d5	8.80	82	327273	27.33	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	286776	39.11	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	974581	26.33	ng	0.00
79) Terphenyl-d14	19.63	244	1493052	26.66	ng	0.00
Target Compounds						
49) Acenaphthylene	14.02	152	233444	4.638	ng	Qvalue 100
71) Phenanthrene	17.09	178	1950590	29.762	ng	99
72) Anthracene	17.18	178	527014	8.402	ng	100
75) Fluoranthene	19.07	202	4104470	54.185	ng	99
78) Pyrene	19.43	202	4362437	54.545	ng	99
81) Benzo(a)anthracene	21.13	228	2448018	31.715	ng	97
83) Chrysene	21.18	228	2300584m	31.129	ng	
87) Indeno(1,2,3-cd)pyrene	25.65	276	1368415	13.980	ng	97
88) Benzo(b)fluoranthene	22.72	252	2629005m	32.226	ng	
89) Benzo(k)fluoranthene	22.75	252	937640m	12.029	ng	
90) Benzo(a)pyrene	23.29	252	2089236	27.605	ng	97
91) Dibenzo(a,h)anthracene	25.65	278	370898	4.613	ng	# 89
92) Benzo(a,h,i)perylene	26.34	276	1343492	16.649	ng	99

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

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