

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090420\
 Data File : BP003211.D
 Acq On : 04 Sep 2020 14:59
 Operator : CG/JU
 Sample : L3877-12 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:47:58 PM

Quant Time: Sep 08 02:44:22 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	7.65	152	310963	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	1226882	20.00	ng	0.00
39) Acenaphthene-d10	14.29	164	741103	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1471949	20.00	ng	0.00
76) Chrysene-d12	21.14	240	1044735	20.00	ng	0.00
86) Perylene-d12	23.38	264	1133100	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	147656	8.40	ng	0.00
7) Phenol-d6	6.84	99	177178	8.04	ng	0.00
23) Nitrobenzene-d5	8.80	82	107815	6.34	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	85230	8.50	ng	0.00
45) 2-Fluorobiphenyl	12.91	172	283539	5.60	ng	0.00
79) Terphenyl-d14	19.62	244	327226	6.89	ng	0.00
Target Compounds						
3) Pyridine	3.59	79	45337	2.562	ng	Qvalue # 44
31) Naphthalene	10.48	128	994152	15.772	ng	99
37) 2-Methylnaphthalene	12.10	142	296132m	6.567	ng	
38) 1-Methylnaphthalene	12.32	142	409485	9.554	ng	99
49) Acenaphthylene	14.01	152	494253	7.183	ng	98
52) Acenaphthene	14.36	154	145997	3.454	ng	99
58) Fluorene	15.35	166	208602	4.079	ng	# 97
71) Phenanthrene	17.09	178	1955288	24.515	ng	100
72) Anthracene	17.18	178	703254	9.213	ng	100
75) Fluoranthene	19.07	202	2563228	27.806	ng	99
78) Pyrene	19.42	202	3966712	58.501	ng	99
81) Benzo(a)anthracene	21.13	228	1958589	29.929	ng	94
83) Chrysene	21.18	228	2181640	34.819	ng	97
87) Indeno(1,2,3-cd)pyrene	25.66	276	1410178	16.688	ng	97
88) Benzo(b)fluoranthene	22.72	252	2221957m	31.550	ng	
89) Benzo(k)fluoranthene	22.74	252	692082m	10.285	ng	
90) Benzo(a)pyrene	23.29	252	2171350	33.234	ng	97
91) Dibenzo(a,h)anthracene	25.66	278	457417	6.590	ng	# 83
92) Benzo(a,h,i)perylene	26.36	276	1819916	26.125	ng	98

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

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