

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092320\
 Data File : BP003364.D
 Acq On : 23 Sep 2020 18:56
 Operator : CG/JU
 Sample : L3872-07DL 4X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BU-02-092120DL

Manual Integrations
 APPROVED

mohammad
 9/24/2020 10:47:12 AM

Quant Time: Sep 24 01:37:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 12:48:08 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	225735	20.00	ng	0.00
21) Naphthalene-d8	10.340	136	902678	20.00	ng	# 0.00
39) Acenaphthene-d10	14.216	164	525651	20.00	ng	0.00
64) Phenanthrene-d10	16.969	188	1006490	20.00	ng	0.00
76) Chrysene-d12	21.069	240	716770	20.00	ng	0.00
86) Perylene-d12	23.263	264	749377	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.181	112	306065	23.68	ng	0.00
7) Phenol-d6	6.763	99	395165	22.93	ng	0.00
23) Nitrobenzene-d5	8.710	82	257634	16.77	ng	-0.01
42) 2,4,6-Tribromophenol	15.716	330	135728	16.93	ng	0.00
45) 2-Fluorobiphenyl	12.828	172	601895	16.75	ng	-0.01
79) Terphenyl-d14	19.545	244	669899	19.47	ng	0.00
Target Compounds						
						Qvalue
49) Acenaphthylene	13.934	152	175871	3.55	ng	99
71) Phenanthrene	17.010	178	1037003	18.94	ng	99
72) Anthracene	17.098	178	434944	8.06	ng	99
75) Fluoranthene	18.998	202	3104990	45.09	ng	100
78) Pyrene	19.351	202	2754982	61.63	ng	100
81) Benzo(a)anthracene	21.051	228	1597456	34.48	ng	98
83) Chrysene	21.104	228	1390773	31.24	ng	98
87) Indeno(1,2,3-cd)pyrene	25.492	276	1160806	20.44	ng	95
88) Benzo(b)fluoranthene	22.610	252	2112467m	44.64	ng	
89) Benzo(k)fluoranthene	22.645	252	718053m	15.83	ng	
90) Benzo(a)pyrene	23.168	252	1576358	35.33	ng	97
91) Dibenzo(a,h)anthracene	25.492	278	295179m	6.30	ng	
92) Benzo(g,h,i)perylene	26.174	276	1111402	23.75	ng	98

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

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