Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070120\
Data File : BP002615.D
Acq On : 01 Jul 2020 21:20
Operator : CG/JU
Sample : L3157-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB-D35

Manual Integrations APPROVED
mohammad
7/7/2020 8:23:37 AM

Quant Time: Jul 02 02:26:15 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 29 16:51:57 2020
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	330802	20.00	ng		0.00
21) Naphthalene-d8	10.34	136	1324590	20.00	ng		0.00
39) Acenaphthene-d10	14.22	164	781835	20.00	ng		0.00
64) Phenanthrene-d10	16.98	188	1528839	20.00	ng		0.00
76) Chrysene-d12	21.09	240	926199	20.00	ng		0.00
86) Perylene-d12	23.29	264	859593	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.19	112	1613659	82.59	ng		0.00
7) Phenol-d6	6.77	99	2245595	82.35	ng		0.00
23) Nitrobenzene-d5	8.72	82	1645677	64.02	ng		0.00
42) 2,4,6-Tribromophenol	15.72	330	731349	76.91	ng		0.00
45) 2-Fluorobiphenyl	12.83	172	3230996	60.82	ng		0.00
79) Terphenyl-d14	19.57	244	3317398	77.31	ng		0.00
Target Compounds							Qvalue
49) Acenaphthylene	13.93	152	177926	2.348	ng		98
50) Dimethylphthalate	13.68	163	620544	10.105	ng		100
70) Pentachlorophenol	16.65	266	768	2.732	ng		97
71) Phenanthrene	17.02	178	907116	11.028	ng		99
72) Anthracene	17.11	178	351347	4.292	ng		100
73) Carbazole	17.39	167	159534	2.017	ng		100
75) Fluoranthene	19.00	202	1315111	13.208	ng		97
78) Pyrene	19.36	202	1262197	20.021	ng		99
81) Benzo(a)anthracene	21.07	228	628568	10.381	ng		94
83) Chrysene	21.12	228	669639m	11.533	ng		
87) Indeno(1,2,3-cd)pyrene	25.50	276	445518	7.198	ng		99
88) Benzo(b)fluoranthene	22.63	252	853376	15.558	ng	#	97
89) Benzo(k)fluoranthene	22.67	252	280062m	5.473	ng		
90) Benzo(a)pyrene	23.19	252	539653	10.721	ng	#	96
91) Dibenzo(a,h)anthracene	25.50	278	119654	2.366	ng	#	76
92) Benzo(g,h,i)perylene	26.17	276	460723	9.109	ng		97

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

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