

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061920\
 Data File : BP002470.D
 Acq On : 19 Jun 2020 21:42
 Operator : CG/JU
 Sample : L3056-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SAMPLE-B-LOT-11

Manual Integrations
 APPROVED

mohammad
 6/22/2020 3:26:23 PM

Quant Time: Jun 20 03:42:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:08:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	211364	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	895794	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	564708	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1222404	20.00	ng	0.00
76) Chrysene-d12	21.10	240	767021	20.00	ng	-0.01
86) Perylene-d12	23.30	264	768924	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1267429	110.96	ng	0.00
7) Phenol-d6	6.78	99	1785403	117.40	ng	0.00
23) Nitrobenzene-d5	8.73	82	1240658	82.71	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	656826	121.48	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2532868	75.42	ng	0.00
79) Terphenyl-d14	19.58	244	3055081	110.81	ng	0.00

Target Compounds					Qvalue	
49) Acenaphthylene	13.95	152	168301	3.483	ng	99
50) Dimethylphthalate	13.69	163	199768	5.163	ng	99
52) Acenaphthene	14.30	154	72114	2.548	ng	99
58) Fluorene	15.29	166	112857	3.148	ng	98
71) Phenanthrene	17.03	178	2099305	36.713	ng	99
72) Anthracene	17.12	178	674614	11.876	ng	98
73) Carbazole	17.39	167	126724	2.268	ng	99
75) Fluoranthene	19.02	202	3498882	51.262	ng	99
78) Pyrene	19.37	202	3122148	66.453	ng	99
81) Benzo(a)anthracene	21.08	228	1489649	34.025	ng	99
83) Chrysene	21.13	228	1347288	32.481	ng	97
87) Indeno(1,2,3-cd)pyrene	25.51	276	811440	16.872	ng	97
88) Benzo(b)fluoranthene	22.64	252	1580891m	38.125	ng	
89) Benzo(k)fluoranthene	22.68	252	549458m	13.946	ng	
90) Benzo(a)pyrene	23.20	252	1187802	30.567	ng	98
91) Dibenzo(a,h)anthracene	25.51	278	209436m	5.429	ng	
92) Benzo(a,h,i)perylene	26.18	276	801392	20.057	ng	98

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

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