

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052920\
 Data File : BP002323.D
 Acq On : 29 May 2020 19:44
 Operator : CG/JU
 Sample : L2746-02 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NMDUP02-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:24:29 PM

Quant Time: May 30 01:56:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 15:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	314366	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1304870	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	778281	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	1628353	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1454741	20.00	ng	0.00
86) Perylene-d12	23.32	264	1520989	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	897604	53.63	ng	0.00
7) Phenol-d6	6.79	99	1227651	53.92	ng	0.00
23) Nitrobenzene-d5	8.75	82	800223	36.87	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	407287	55.22	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	1751289	39.71	ng	0.00
79) Terphenyl-d14	19.59	244	2476603	41.23	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	202358	3.841	ng	99
71) Phenanthrene	17.05	178	796893	10.332	ng	99
72) Anthracene	17.14	178	242558	3.196	ng	99
75) Fluoranthene	19.03	202	1176097	13.239	ng	99
78) Pyrene	19.38	202	1100534	12.149	ng	98
81) Benzo(a)anthracene	21.09	228	567249	6.812	ng	100
83) Chrysene	21.15	228	503793	6.304	ng	97
87) Indeno(1,2,3-cd)pyrene	25.53	276	308934	3.281	ng	97
88) Benzo(b)fluoranthene	22.66	252	652961m	7.874	ng	
89) Benzo(k)fluoranthene	22.70	252	207060m	2.602	ng	
90) Benzo(a)pyrene	23.22	252	491352	6.388	ng	# 94
92) Benzo(a,h,i)perylene	26.21	276	304989	3.895	ng	# 93

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

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