

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052820\
 Data File : BP002303.D
 Acq On : 28 May 2020 11:25
 Operator : CG/JU
 Sample : L2744-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM55-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 5/29/2020 10:30:51 PM

Quant Time: May 28 12:38:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 28 11:01:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	296606	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1232747	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	718417	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	1491480	20.00	ng	0.00
76) Chrysene-d12	21.09	240	1377312	20.00	ng	0.00
86) Perylene-d12	23.29	264	1419622	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1591340	100.76	ng	0.00
7) Phenol-d6	6.79	99	2214113	103.06	ng	0.00
23) Nitrobenzene-d5	8.75	82	1451806	70.81	ng	0.00
42) 2,4,6-Tribromophenol	15.74	330	676529	99.36	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	2992519	73.50	ng	0.00
79) Terphenyl-d14	19.59	244	3900220	68.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	564607	11.611	ng	Qvalue 100
71) Phenanthrene	17.04	178	212850	3.013	ng	98
75) Fluoranthene	19.02	202	580316	7.132	ng	99
78) Pyrene	19.37	202	570350	6.650	ng	98
81) Benzo(a)anthracene	21.08	228	352255	4.468	ng	98
83) Chrysene	21.13	228	394784	5.217	ng	97
87) Indeno(1,2,3-cd)pyrene	25.50	276	298022	3.392	ng	96
88) Benzo(b)fluoranthene	22.64	252	682399m	8.817	ng	
89) Benzo(k)fluoranthene	22.67	252	228465m	3.076	ng	
90) Benzo(a)pyrene	23.19	252	464535	6.470	ng	# 97
92) Benzo(g,h,i)perylene	26.17	276	290492	3.975	ng	# 93

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

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