

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052820\
 Data File : BP002312.D
 Acq On : 28 May 2020 16:33
 Operator : CG/JU
 Sample : L2744-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM78-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 5/29/2020 10:31:13 PM

Quant Time: May 29 03:13:35 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	285596	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	1173302	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	696001	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	1464750	20.00	ng	0.00
76) Chrysene-d12	21.10	240	1351745	20.00	ng	0.00
86) Perylene-d12	23.30	264	1419383	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1207218	79.39	ng	0.00
7) Phenol-d6	6.79	99	1666005	80.54	ng	0.00
23) Nitrobenzene-d5	8.75	82	1074649	55.07	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	532515	80.73	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	2269509	57.54	ng	0.00
79) Terphenyl-d14	19.59	244	3110042	55.72	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	427345	9.071	ng	Qvalue 100
71) Phenanthrene	17.04	178	234598	3.381	ng	98
75) Fluoranthene	19.02	202	621646	7.779	ng	99
78) Pyrene	19.37	202	596749	7.090	ng	99
81) Benzo(a)anthracene	21.08	228	351477	4.542	ng	98
83) Chrysene	21.13	228	328265	4.420	ng	97
87) Indeno(1,2,3-cd)pyrene	25.50	276	240661	2.739	ng	96
88) Benzo(b)fluoranthene	22.64	252	482644	6.237	ng	# 97
89) Benzo(k)fluoranthene	22.67	252	150335m	2.024	ng	
90) Benzo(a)pyrene	23.20	252	343974	4.792	ng	# 97
92) Benzo(g,h,i)perylene	26.17	276	240513	3.292	ng	# 91

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

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