

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002340.D
 Acq On : 01 Jun 2020 15:44
 Operator : CG/JU
 Sample : L2744-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM79-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:47 PM

Quant Time: Jun 01 16:17:37 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	7.61	152	296751	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1222944	20.00	ng	0.00
39) Acenaphthene-d10	14.26	164	734155	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1561323	20.00	ng	0.00
76) Chrysene-d12	21.12	240	1340462	20.00	ng	0.00
86) Perylene-d12	23.32	264	1396377	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	1763391	111.60	ng	0.00
7) Phenol-d6	6.80	99	2405929	111.94	ng	0.00
23) Nitrobenzene-d5	8.75	82	1595267	78.43	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	800982	115.12	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	3298302	79.27	ng	0.00
79) Terphenyl-d14	19.60	244	4171538	75.36	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.72	163	542936	10.926	ng	99
71) Phenanthrene	17.05	178	286094	3.868	ng	99
75) Fluoranthene	19.04	202	732711	8.602	ng	99
78) Pyrene	19.39	202	722560	8.657	ng	98
81) Benzo(a)anthracene	21.10	228	391305	5.100	ng	98
83) Chrysene	21.15	228	358520	4.868	ng	97
87) Indeno(1,2,3-cd)pyrene	25.54	276	252864	2.926	ng	98
88) Benzo(b)fluoranthene	22.66	252	511015m	6.712	ng	
89) Benzo(k)fluoranthene	22.70	252	171212m	2.344	ng	
90) Benzo(a)pyrene	23.23	252	369800	5.237	ng	# 96
92) Benzo(g,h,i)perylene	26.22	276	253912	3.532	ng	# 88

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

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