

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002338.D
 Acq On : 01 Jun 2020 14:36
 Operator : CG/JU
 Sample : L2773-21
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM91-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:36 PM

Quant Time: Jun 01 15:20:36 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	392730	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1599145	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	948584	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	2004528	20.00	ng	0.00
76) Chrysene-d12	21.12	240	1721883	20.00	ng	0.00
86) Perylene-d12	23.33	264	1825040	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	1810801	86.60	ng	0.00
7) Phenol-d6	6.80	99	2509638	88.23	ng	0.00
23) Nitrobenzene-d5	8.75	82	1687095	63.43	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	835918	92.98	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	3443700	64.06	ng	0.00
79) Terphenyl-d14	19.60	244	4263757	59.97	ng	0.00
Target Compounds						
32) Benzoic acid	9.65	122	3143	4.989	ng	87
50) Dimethylphthalate	13.72	163	449052	6.994	ng	99
71) Phenanthrene	17.05	178	828261	8.723	ng	98
72) Anthracene	17.14	178	228995	2.451	ng	98
75) Fluoranthene	19.04	202	1630160	14.906	ng	100
78) Pyrene	19.39	202	1485686	13.856	ng	98
81) Benzo(a)anthracene	21.10	228	749373	7.603	ng	99
83) Chrysene	21.15	228	690539	7.300	ng	98
87) Indeno(1,2,3-cd)pyrene	25.54	276	481960	4.266	ng	98
88) Benzo(b)fluoranthene	22.66	252	1015827m	10.209	ng	
89) Benzo(k)fluoranthene	22.70	252	253047m	2.650	ng	
90) Benzo(a)pyrene	23.23	252	708324	7.674	ng	97
92) Benzo(a,h,i)perylene	26.23	276	499702	5.319	ng	# 92

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

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