

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002337.D
 Acq On : 01 Jun 2020 14:02
 Operator : CG/JU
 Sample : L2773-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM33-COMP-2020-0-6

Quant Time: Jun 02 06:52:01 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	407509	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1670267	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	988350	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	2053252	20.00	ng	0.00
76) Chrysene-d12	21.12	240	1821924	20.00	ng	0.00
86) Perylene-d12	23.33	264	1904914	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.23	112	2334785	107.61	ng	0.00
7) Phenol-d6	6.80	99	3203591	108.54	ng	0.00
23) Nitrobenzene-d5	8.75	82	2190808	78.86	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	1076885	114.96	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	4290615	76.60	ng	0.00
79) Terphenyl-d14	19.60	244	5237471	69.62	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
32) Benzoic acid	9.65	122	1948	4.916	ng	92
50) Dimethylphthalate	13.72	163	462871	6.919	ng	100
71) Phenanthrene	17.05	178	458213	4.711	ng	99
75) Fluoranthene	19.03	202	919417	8.208	ng	99
78) Pyrene	19.39	202	868406	7.654	ng	99
80) Butylbenzylphthalate	20.28	149	211363	4.679	ng	92
81) Benzo(a)anthracene	21.10	228	439826	4.217	ng	95
83) Chrysene	21.15	228	398486	3.981	ng	98
87) Indeno(1,2,3-cd)pyrene	25.54	276	281653	2.389	ng	97
88) Benzo(b)fluoranthene	22.66	252	569795	5.486	ng	# 96
90) Benzo(a)pyrene	23.23	252	416057	4.319	ng	# 97
92) Benzo(a,h,i)perylene	26.22	276	294652	3.005	ng	# 94

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

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