

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072220\
 Data File : BP002829.D
 Acq On : 22 Jul 2020 17:32
 Operator : CG/JU
 Sample : L3362-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-59

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:11:03 PM

Quant Time: Jul 22 18:53:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	122948	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	470999	20.00	ng	-0.01
39) Acenaphthene-d10	14.20	164	230768	20.00	ng	-0.01
64) Phenanthrene-d10	16.98	188	200553	20.00	ng	0.01
76) Chrysene-d12	21.11	240	175652	20.00	ng	0.04
86) Perylene-d12	23.30	264	298462	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	566579	77.75	ng	0.00
7) Phenol-d6	6.75	99	730399	70.23	ng	0.00
23) Nitrobenzene-d5	8.69	82	484488	55.02	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	139243	57.80	ng	0.00
45) 2-Fluorobiphenyl	12.81	172	937836	62.87	ng	-0.01
79) Terphenyl-d14	19.59	244	459022	59.91	ng	0.03
Target Compounds						
32) Benzoic acid	9.69	122	10414	8.779	ng	94
50) Dimethylphthalate	13.66	163	190414	10.819	ng	# 94
71) Phenanthrene	17.02	178	56519	5.205	ng	# 15
75) Fluoranthene	19.03	202	91387	7.263	ng	# 55
78) Pyrene	19.39	202	87953	7.252	ng	# 58
81) Benzo(a)anthracene	21.10	228	46548	4.094	ng	# 60
83) Chrysene	21.14	228	56615	5.231	ng	# 59
87) Indeno(1,2,3-cd)pyrene	25.50	276	90969	4.217	ng	98
88) Benzo(b)fluoranthene	22.65	252	114891m	6.199	ng	
90) Benzo(a)pyrene	23.21	252	74672	4.252	ng	# 59
92) Benzo(g,h,i)perylene	26.17	276	93047	5.275	ng	# 89

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

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