

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072020\
 Data File : BP002797.D
 Acq On : 20 Jul 2020 13:13
 Operator : CG/JU
 Sample : L3177-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AU-05-071720

Manual Integrations
 APPROVED

mohammad
 7/21/2020 11:33:05 AM

Quant Time: Jul 20 15:31:26 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	167939	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	682716	20.00	ng	-0.01
39) Acenaphthene-d10	14.20	164	410203	20.00	ng	0.00
64) Phenanthrene-d10	16.96	188	836080	20.00	ng	0.00
76) Chrysene-d12	21.07	240	722909	20.00	ng	0.00
86) Perylene-d12	23.27	264	600604	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	975081	97.96	ng	0.00
7) Phenol-d6	6.76	99	1335024	93.98	ng	0.00
23) Nitrobenzene-d5	8.69	82	837105	65.58	ng	-0.01
42) 2,4,6-Tribromophenol	15.71	330	424594	99.16	ng	0.00
45) 2-Fluorobiphenyl	12.82	172	1649457	62.21	ng	0.00
79) Terphenyl-d14	19.56	244	1918667	60.85	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.66	163	208738	6.672	ng	99
71) Phenanthrene	17.00	178	98507	2.176	ng	99
75) Fluoranthene	19.00	202	190281	3.628	ng	99
78) Pyrene	19.35	202	221448	4.437	ng	97
81) Benzo(a)anthracene	21.06	228	97873	2.091	ng	# 93
83) Chrysene	21.11	228	100729m	2.261	ng	
88) Benzo(b)fluoranthene	22.62	252	104616	2.805	ng	# 73
90) Benzo(a)pyrene	23.17	252	83985	2.376	ng	# 86
92) Benzo(g,h,i)perylene	26.15	276	71605	2.017	ng	92

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

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