

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031320\
 Data File : BG044770.D
 Acq On : 13 Mar 2020 16:56
 Operator : CG/JU
 Sample : L1755-01 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-03-031020

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:51:46 PM

Quant Time: Mar 13 17:38:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 16:27:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	146370	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	596086	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	387556	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	775035	20.00	ng	0.00
76) Chrysene-d12	21.70	240	630240	20.00	ng	-0.01
87) Perylene-d12	24.93	264	710865	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	353423	37.59	ng	0.00
7) Phenol-d6	7.21	99	534754	39.14	ng	0.00
23) Nitrobenzene-d5	9.20	82	346136	25.48	ng	-0.01
42) 2,4,6-Tribromophenol	16.17	330	196427	38.71	ng	0.00
45) 2-Fluorobiphenyl	13.31	172	755060	27.90	ng	0.00
79) Terphenyl-d14	20.04	244	1013644	30.08	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.13	163	85739	2.778	ng	96
75) Fluoranthene	19.48	202	152153	3.085	ng	91
78) Pyrene	19.84	202	166647	3.604	ng	99
81) Benzo(a)anthracene	21.68	228	94125	2.074	ng	92
83) Chrysene	21.75	228	110478m	2.549	ng	
88) Benzo(b)fluoranthene	23.91	252	145071	3.091	ng	# 95
90) Benzo(a)pyrene	24.77	252	111058	2.529	ng	# 94

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

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