

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052419\
 Data File : BM020562.D
 Acq On : 25 May 2019 03:56
 Operator : JU/SJ
 Sample : K2647-11 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SU-03-052219-A

Manual Integrations
 APPROVED

mohammad
 5/27/2019 4:00:26 AM

Quant Time: May 25 06:38:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 24 16:07:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	70269	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	264095	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	201493	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	563887	20.00	ng	0.00
76) Chrysene-d12	21.26	240	531646	20.00	ng	-0.01
87) Perylene-d12	23.50	264	539503	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	213113	59.50	ng	0.00
7) Phenol-d6	6.84	99	294429	54.46	ng	0.00
23) Nitrobenzene-d5	8.83	82	189118	30.06	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	215043	53.27	ng	0.00
45) 2-Fluorobiphenyl	12.93	172	522815	32.56	ng	0.00
79) Terphenyl-d14	19.70	244	1169404	39.76	ng	-0.01
Target Compounds						
49) Acenaphthylene	14.04	152	42133	2.106	ng	98
50) Dimethylphthalate	13.77	163	58106	3.256	ng	98
71) Phenanthrene	17.11	178	227737	7.434	ng	97
72) Anthracene	17.20	178	75083	2.527	ng	100
75) Fluoranthene	19.13	202	648276	15.218	ng	100
78) Pyrene	19.50	202	634585	19.013	ng	99
81) Benzo(a)anthracene	21.24	228	344989	9.376	ng	96
83) Chrysene	21.30	228	324988	9.140	ng	96
86) Indeno(1,2,3-cd)pyrene	25.77	276	208718	4.855	ng	96
88) Benzo(b)fluoranthene	22.83	252	435111m	12.322	ng	
89) Benzo(k)fluoranthene	22.86	252	144044m	4.203	ng	
90) Benzo(a)pyrene	23.40	252	310326	9.550	ng	# 95
92) Benzo(a,h,i)perylene	26.46	276	217041	7.039	ng	100

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

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