

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050619\
 Data File : BM020152.D
 Acq On : 07 May 2019 04:03
 Operator : JU/SJ
 Sample : K2673-12 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-16-COMP

Manual Integrations
 APPROVED

mohammad
 5/7/2019 3:12:00 PM

Quant Time: May 07 05:38:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	181688	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	644403	20.00	ng	0.00
39) Acenaphthene-d10	14.43	164	348431	20.00	ng	0.00
64) Phenanthrene-d10	17.18	188	791972	20.00	ng	0.00
76) Chrysene-d12	21.37	240	800795	20.00	ng	0.00
87) Perylene-d12	23.66	264	943396	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	271119	24.39	ng	0.00
7) Phenol-d6	6.96	99	367301	24.19	ng	0.00
23) Nitrobenzene-d5	8.96	82	242941	14.88	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	123422	22.82	ng	0.00
45) 2-Fluorobiphenyl	13.05	172	415855	14.68	ng	0.00
79) Terphenyl-d14	19.80	244	577403	12.58	ng	0.00
Target Compounds						
49) Acenaphthylene	14.16	152	150367	4.103	ng	99
50) Dimethylphthalate	13.89	163	83063	2.805	ng	97
71) Phenanthrene	17.23	178	682603	15.614	ng	99
72) Anthracene	17.32	178	220569	5.046	ng	99
75) Fluoranthene	19.24	202	1328390	24.016	ng	99
78) Pyrene	19.60	202	1144494	21.287	ng	100
81) Benzo(a)anthracene	21.35	228	783149	13.945	ng	97
83) Chrysene	21.40	228	707518	12.990	ng	97
86) Indeno(1,2,3-cd)pyrene	26.00	276	593590	9.162	ng	# 100
88) Benzo(b)fluoranthene	22.97	252	1118406m	17.953	ng	
89) Benzo(k)fluoranthene	23.01	252	392172m	6.901	ng	
90) Benzo(a)pyrene	23.56	252	749794	13.251	ng	# 93
91) Dibenzo(a,h)anthracene	26.01	278	155883	2.649	ng	# 79
92) Benzo(a,h,i)perylene	26.72	276	530663	9.498	ng	95

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

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