

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
 Data File : BM020115.D
 Acq On : 03 May 2019 17:04
 Operator : JU/SJ
 Sample : K2584-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSSI-0419-S-DH-8

Quant Time: May 03 17:42:48 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	146195	20.00	ng	0.00
21) Naphthalene-d8	10.58	136	527842	20.00	ng	0.00
39) Acenaphthene-d10	14.43	164	316931	20.00	ng	0.00
64) Phenanthrene-d10	17.17	188	733993	20.00	ng	0.00
76) Chrysene-d12	21.36	240	777487	20.00	ng	-0.01
87) Perylene-d12	23.64	264	924100	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.36	112	1137767	127.21	ng	0.00
7) Phenol-d6	6.95	99	1745348	142.84	ng	0.00
23) Nitrobenzene-d5	8.95	82	1266388	94.72	ng	0.00
42) 2,4,6-Tribromophenol	15.92	330	804318	163.50	ng	0.00
45) 2-Fluorobiphenyl	13.05	172	2597219	100.83	ng	-0.01
79) Terphenyl-d14	19.80	244	4593875	103.06	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
50) Dimethylphthalate	13.88	163	165510	6.144	ng	99
71) Phenanthrene	17.22	178	314304	7.757	ng	99
75) Fluoranthene	19.23	202	481858	9.399	ng	99
78) Pyrene	19.60	202	409989	7.854	ng	100
81) Benzo(a)anthracene	21.34	228	254490	4.667	ng	97
83) Chrysene	21.39	228	221442	4.188	ng	95
86) Indeno(1,2,3-cd)pyrene	25.96	276	128536	2.044	ng	# 100
88) Benzo(b)fluoranthene	22.95	252	289953	4.752	ng	# 93
90) Benzo(a)pyrene	23.54	252	200165	3.611	ng	# 93
92) Benzo(g,h,i)perylene	26.67	276	116952	2.137	ng	# 96

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

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