

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082119\
 Data File : BM022230.D
 Acq On : 21 Aug 2019 15:43
 Operator : HP/JU
 Sample : K4421-13
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SUNSET-PARK-SB-1-COMP

Quant Time: Aug 21 16:19:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 20 02:18:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	33197	20.00	ng	-0.01
21) Naphthalene-d8	10.34	136	128941	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	76287	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	146114	20.00	ng	0.00
76) Chrysene-d12	21.19	240	147899	20.00	ng	0.00
87) Perylene-d12	23.39	264	173063	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	183314	98.40	ng	0.00
7) Phenol-d6	6.76	99	245150	98.82	ng	0.00
23) Nitrobenzene-d5	8.73	82	173241	65.35	ng	-0.01
42) 2,4,6-Tribromophenol	15.73	330	105987	85.84	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	398142	65.19	ng	-0.01
79) Terphenyl-d14	19.63	244	614913	60.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.70	163	55857	8.709	ng	98
71) Phenanthrene	17.03	178	47249	5.576	ng	97
75) Fluoranthene	19.06	202	60861	5.956	ng	98
78) Pyrene	19.42	202	56916	5.467	ng	98
81) Benzo(a)anthracene	21.17	228	31543	3.033	ng	99
83) Chrysene	21.23	228	29556	2.944	ng	96
84) Bis(2-ethylhexyl)phthalate	21.10	149	15637	2.233	ng	# 95
86) Indeno(1,2,3-cd)pyrene	25.58	276	22831	2.030	ng	# 94
88) Benzo(b)fluoranthene	22.73	252	40365	3.493	ng	98
90) Benzo(a)pyrene	23.29	252	28322	2.691	ng	# 96
92) Benzo(g,h,i)perylene	26.26	276	22595	2.311	ng	# 96

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

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