

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM083019\
 Data File : BM022460.D
 Acq On : 31 Aug 2019 02:50
 Operator : HP/JU
 Sample : K4439-04 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 S-1(5-7)

Quant Time: Aug 31 05:54:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 16:51:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	20766	20.00	ng	-0.01
21) Naphthalene-d8	10.29	136	74125	20.00	ng	0.00
39) Acenaphthene-d10	14.18	164	46777	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	107972	20.00	ng	0.00
76) Chrysene-d12	21.16	240	171683	20.00	ng	-0.01
87) Perylene-d12	23.36	264	220065	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	43692	36.93	ng	0.00
7) Phenol-d6	6.72	99	55315	34.96	ng	0.00
23) Nitrobenzene-d5	8.70	82	37980	22.44	ng	0.00
42) 2,4,6-Tribromophenol	15.70	330	27284	28.50	ng	0.00
45) 2-Fluorobiphenyl	12.79	172	84522	20.47	ng	0.00
79) Terphenyl-d14	19.59	244	189024	14.53	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.66	163	11873	2.918	ng	Qvalue # 94
71) Phenanthrene	16.99	178	15016	2.386	ng	97
75) Fluoranthene	19.02	202	29825	3.559	ng	97
78) Pyrene	19.39	202	31952	2.731	ng	97
81) Benzo(a)anthracene	21.15	228	25424	2.063	ng	# 88
83) Chrysene	21.20	228	33588	2.864	ng	# 90
88) Benzo(b)fluoranthene	22.70	252	39498	2.584	ng	# 86

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

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