

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053019\
 Data File : BF114684.D
 Acq On : 31 May 2019 2:05
 Operator : HP/JU
 Sample : K3041-11DL 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-DDL

Quant Time: May 31 07:52:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	386693	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	1427675	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	633260	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	1132392	20.00	ng	0.00
76) Chrysene-d12	13.82	240	744055	20.00	ng	0.00
87) Perylene-d12	15.21	264	613901	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1359498	61.73	ng	0.00
7) Phenol-d6	6.33	99	1640765	61.82	ng	0.00
23) Nitrobenzene-d5	7.25	82	960999	42.00	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	324123	53.72	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1432258	32.63	ng	0.00
79) Terphenyl-d14	12.78	244	1252902	31.66	ng	0.00

Target Compounds				Qvalue		
20) 3+4-Methylphenols	7.10	107	52575	2.078	ng	# 56
31) Naphthalene	8.00	128	1193557	18.150	ng	97
37) 2-Methylnaphthalene	8.69	142	516611	11.824	ng	98
38) 1-Methylnaphthalene	8.79	142	238767	5.768	ng	99
50) Dimethylphthalate	9.45	163	199458	4.520	ng	97
52) Acenaphthene	9.76	154	800475	21.911	ng	99
55) Dibenzofuran	9.93	168	976838	18.784	ng	97
58) Fluorene	10.27	166	1144426	30.138	ng	100
71) Phenanthrene	11.23	178	3780782	65.451	ng	98
72) Anthracene	11.27	178	1196983	20.474	ng	98
73) Carbazole	11.42	167	283490	5.517	ng	95
75) Fluoranthene	12.40	202	1684002	28.154	ng	98
78) Pyrene	12.63	202	1117783	17.759	ng	98
81) Benzo(a)anthracene	13.81	228	182131	3.185	ng	98
83) Chrysene	13.85	228	159691	2.923	ng	93

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

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