

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113469.D
 Acq On : 1 Apr 2019 16:28
 Operator : JU/SJ
 Sample : K2056-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-10

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:49:22 PM

Quant Time: Apr 02 07:12:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	136729	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	508934	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	241069	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	463467	20.00	ng	0.00
76) Chrysene-d12	13.84	240	355680	20.00	ng	0.00
87) Perylene-d12	15.25	264	408716	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	935230	111.84	ng	0.00
7) Phenol-d6	6.36	99	1203402	113.76	ng	0.00
23) Nitrobenzene-d5	7.27	82	732675	85.45	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	322853	108.42	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1371964	90.30	ng	0.00
79) Terphenyl-d14	12.79	244	1332445	82.57	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.46	163	94253	5.276	ng	Qvalue 100
71) Phenanthrene	11.24	178	207608	9.020	ng	97
72) Anthracene	11.29	178	72282	3.092	ng	98
73) Carbazole	11.45	167	82164	3.683	ng	98
75) Fluoranthene	12.43	202	1022357	39.293	ng	98
78) Pyrene	12.65	202	745554	30.249	ng	99
81) Benzo(a)anthracene	13.83	228	126049	6.191	ng	95
83) Chrysene	13.87	228	385694	18.651	ng	99
86) Indeno(1,2,3-cd)pyrene	16.61	276	104503	5.888	ng	96
88) Benzo(b)fluoranthene	14.85	252	419163m	16.509	ng	
89) Benzo(k)fluoranthene	14.87	252	126655m	5.680	ng	
90) Benzo(a)pyrene	15.19	252	81303	3.612	ng	97
92) Benzo(g,h,i)perylene	17.02	276	88482	4.509	ng	96

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

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