

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032719\
 Data File : BF113341.D
 Acq On : 27 Mar 2019 20:07
 Operator : JU/SJ
 Sample : K2097-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-6

Manual Integrations
 APPROVED

Sohil
 3/28/2019 8:41:20 AM

Quant Time: Mar 28 04:16:47 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	180772	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	678743	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	318033	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	592424	20.00	ng	0.00
76) Chrysene-d12	13.83	240	426642	20.00	ng	-0.01
87) Perylene-d12	15.24	264	438347	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	956182	86.49	ng	0.00
7) Phenol-d6	6.35	99	1238175	88.53	ng	0.00
23) Nitrobenzene-d5	7.26	82	763691	66.78	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	332780	84.71	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1424841	66.35	ng	0.00
79) Terphenyl-d14	12.79	244	1284301	66.35	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.45	163	98019	4.159	ng	99
71) Phenanthrene	11.23	178	431062	14.652	ng	98
72) Anthracene	11.29	178	80645	2.699	ng	100
73) Carbazole	11.44	167	85452	2.997	ng	98
75) Fluoranthene	12.42	202	1152218	34.644	ng	99
78) Pyrene	12.64	202	874109	29.566	ng	99
81) Benzo(a)anthracene	13.82	228	138689	5.679	ng	94
83) Chrysene	13.86	228	486002	19.592	ng	99
86) Indeno(1,2,3-cd)pyrene	16.58	276	125668	5.903	ng	98
88) Benzo(b)fluoranthene	14.84	252	506767m	18.611	ng	
89) Benzo(k)fluoranthene	14.87	252	182095m	7.615	ng	
90) Benzo(a)pyrene	15.18	252	114534	4.744	ng	100
92) Benzo(a,h,i)perylene	16.99	276	108000	5.132	ng	96

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

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