

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041519\
 Data File : BF113775.D
 Acq On : 15 Apr 2019 21:03
 Operator : JU/SJ
 Sample : K2397-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-S

Manual Integrations
 APPROVED

Sohil
 4/16/2019 5:02:59 PM

Quant Time: Apr 16 00:46:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	115349	20.00	ng	-0.02
21) Naphthalene-d8	7.95	136	428201	20.00	ng	-0.02
39) Acenaphthene-d10	9.70	164	203338	20.00	ng	-0.02
64) Phenanthrene-d10	11.19	188	377523	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	258335	20.00	ng	-0.02
87) Perylene-d12	15.24	264	338307	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	742570	109.78	ng	-0.01
7) Phenol-d6	6.33	99	983508	117.97	ng	-0.01
23) Nitrobenzene-d5	7.24	82	584003	79.89	ng	-0.02
42) 2,4,6-Tribromophenol	10.49	330	256244	104.91	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	1050593	83.18	ng	-0.02
79) Terphenyl-d14	12.77	244	878249	69.22	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	7.97	128	47413	2.285	ng	99
50) Dimethylphthalate	9.43	163	102670	6.850	ng	99
71) Phenanthrene	11.21	178	366408	19.532	ng	99
72) Anthracene	11.27	178	109175	5.720	ng	98
75) Fluoranthene	12.40	202	886381	44.094	ng	99
78) Pyrene	12.63	202	878497	44.693	ng	100
81) Benzo(a)anthracene	13.82	228	688703	46.215	ng	99
83) Chrysene	13.85	228	591488	39.154	ng	98
86) Indeno(1,2,3-cd)pyrene	16.61	276	587216	42.101	ng	96
88) Benzo(b)fluoranthene	14.84	252	998113m	48.198	ng	
89) Benzo(k)fluoranthene	14.86	252	356675m	19.191	ng	
90) Benzo(a)pyrene	15.19	252	800852	43.521	ng	98
91) Dibenzo(a,h)anthracene	16.60	278	148956m	8.656	ng	
92) Benzo(g,h,i)perylene	17.02	276	570366	32.473	ng	97

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

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