

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041619\
 Data File : BF113795.D
 Acq On : 16 Apr 2019 15:57
 Operator : JU/SJ
 Sample : K2411-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-14-S

Manual Integrations
 APPROVED

Sohil
 4/17/2019 3:45:35 PM

Quant Time: Apr 16 18:18:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 16 17:22:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	3968	20.00	ng	0.00
21) Naphthalene-d8	0.00	136	0	0.00	ng	-7.96
39) Acenaphthene-d10	9.43	164	8395	20.00	ng	-0.28
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-11.19
76) Chrysene-d12	13.83	240	273m	20.00	ng	0.00
87) Perylene-d12	15.24	264	116	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	618297	2657.10	ng	0.00
7) Phenol-d6	6.33	99	815105	2842.08	ng	0.00
23) Nitrobenzene-d5	7.24	82	457019	0.00	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	155856m	1545.51	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	784212	1503.87	ng	0.00
79) Terphenyl-d14	12.77	244	726035	54147.20	ng	0.00
Target Compounds						
15) Benzyl Alcohol	6.83	79	7140	36.819	ng	Qvalue # 40
50) Dimethylphthalate	9.43	163	92899m	150.122	ng	
78) Pyrene	12.63	202	41609	2003.104	ng	98
81) Benzo(a)anthracene	13.82	228	31381	1992.682	ng	95
83) Chrysene	13.86	228	24727	1548.908	ng	97
84) Bis(2-ethylhexyl)phthalate	13.80	149	3441	336.587	ng	# 89
86) Indeno(1,2,3-cd)pyrene	16.63	276	24521	1663.632	ng	99
88) Benzo(b)fluoranthene	14.85	252	43503m	6126.634	ng	
89) Benzo(k)fluoranthene	14.86	252	18683m	2931.745	ng	
90) Benzo(a)pyrene	15.19	252	35948	5697.326	ng	95
91) Dibenzo(a,h)anthracene	16.61	278	7206	1221.263	ng	# 74
92) Benzo(g,h,i)perylene	17.04	276	21052	3495.544	ng	94

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

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