

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051219\
 Data File : BF114199.D
 Acq On : 12 May 2019 18:14
 Operator : JU/SJ
 Sample : K2767-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS010-0206-01

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:37:32 PM

Quant Time: May 13 09:12:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	309779	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	1157066	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	532899	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	977248	20.00	ng	0.00
76) Chrysene-d12	14.02	240	608423	20.00	ng	0.00
87) Perylene-d12	15.49	264	715604	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1398668	72.66	ng	0.00
7) Phenol-d6	6.50	99	1849460	81.23	ng	0.00
23) Nitrobenzene-d5	7.42	82	1154280	55.98	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	412974	74.96	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1751461	49.72	ng	0.00
79) Terphenyl-d14	12.96	244	1330326	45.07	ng	0.00
Target Compounds						
31) Naphthalene	8.15	128	166865	3.087	ng	97
32) Benzoic acid	7.86	122	5174m	7.408	ng	
37) 2-Methylnaphthalene	8.85	142	72882	2.090	ng	94
49) Acenaphthylene	9.75	152	376477	7.144	ng	98
50) Dimethylphthalate	9.60	163	424107	10.841	ng	100
71) Phenanthrene	11.40	178	1125766	23.678	ng	98
72) Anthracene	11.45	178	199736	4.183	ng	96
75) Fluoranthene	12.59	202	1634916	31.932	ng	97
78) Pyrene	12.82	202	2714177	55.454	ng	99
81) Benzo(a)anthracene	14.00	228	1049349m	26.665	ng	
83) Chrysene	14.04	228	1232427	31.948	ng	99
86) Indeno(1,2,3-cd)pyrene	16.96	276	646687	18.968	ng	97
88) Benzo(b)fluoranthene	15.06	252	1558766m	34.000	ng	
89) Benzo(k)fluoranthene	15.09	252	394986m	9.379	ng	
90) Benzo(a)pyrene	15.43	252	947430	23.481	ng	100
91) Dibenzo(a,h)anthracene	16.96	278	186647m	5.567	ng	
92) Benzo(g,h,i)perylene	17.41	276	700382	20.845	ng	97

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

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