

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
 Data File : BF114165.D
 Acq On : 11 May 2019 20:55
 Operator : JU/SJ
 Sample : K2765-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS004-1824-01

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:36:26 PM

Quant Time: May 13 08:01:26 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	251287	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	952926	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	450658	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	916371	20.00	ng	0.00
76) Chrysene-d12	14.00	240	712753	20.00	ng	-0.02
87) Perylene-d12	15.47	264	741200	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1775956	113.73	ng	0.00
7) Phenol-d6	6.49	99	2386121	129.20	ng	0.00
23) Nitrobenzene-d5	7.41	82	1449906	85.39	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	573691	123.13	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2206913	74.08	ng	0.00
79) Terphenyl-d14	12.95	244	2358901	68.22	ng	0.00
Target Compounds						
49) Acenaphthylene	9.74	152	95315	2.139	ng	98
50) Dimethylphthalate	9.59	163	381139	11.521	ng	99
71) Phenanthrene	11.39	178	249931	5.606	ng	98
75) Fluoranthene	12.58	202	519094	10.812	ng	99
78) Pyrene	12.80	202	934267	16.294	ng	98
81) Benzo(a)anthracene	13.99	228	361726m	7.846	ng	
83) Chrysene	14.03	228	445468	9.857	ng	96
86) Indeno(1,2,3-cd)pyrene	16.93	276	299616	7.502	ng	97
88) Benzo(b)fluoranthene	15.04	252	485977m	10.234	ng	
89) Benzo(k)fluoranthene	15.07	252	152070m	3.486	ng	
90) Benzo(a)pyrene	15.41	252	299159	7.158	ng	99
91) Dibenzo(a,h)anthracene	16.93	278	77793m	2.240	ng	
92) Benzo(a,h,i)perylene	17.37	276	360287	10.353	ng	97

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

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