

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
 Data File : BF114167.D
 Acq On : 11 May 2019 21:48
 Operator : JU/SJ
 Sample : K2765-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS002-1218-02

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:35:55 PM

Quant Time: May 13 08:05:38 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	271336	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	957791	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	462130	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	948420	20.00	ng	0.00
76) Chrysene-d12	14.00	240	664015	20.00	ng	-0.02
87) Perylene-d12	15.47	264	727216	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	1975326	117.15	ng	0.00
7) Phenol-d6	6.49	99	2638398	132.30	ng	0.00
23) Nitrobenzene-d5	7.41	82	1541395	90.32	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	626839	131.20	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2403919	78.69	ng	0.00
79) Terphenyl-d14	12.95	244	2374640	73.72	ng	0.00

Target Compounds

						Qvalue
49) Acenaphthylene	9.74	152	116362	2.546	ng	97
50) Dimethylphthalate	9.59	163	524308	15.455	ng	99
71) Phenanthrene	11.39	178	740845	16.056	ng	98
75) Fluoranthene	12.58	202	1015213	20.431	ng	99
78) Pyrene	12.81	202	1799013	33.679	ng	99
81) Benzo(a)anthracene	13.99	228	529589m	12.331	ng	
83) Chrysene	14.03	228	607353	14.426	ng	98
86) Indeno(1,2,3-cd)pyrene	16.92	276	314676	8.457	ng	95
88) Benzo(b)fluoranthene	15.04	252	604920m	12.984	ng	
89) Benzo(k)fluoranthene	15.07	252	157613m	3.683	ng	
90) Benzo(a)pyrene	15.41	252	375395	9.155	ng	98
91) Dibenzo(a,h)anthracene	16.93	278	79313	2.328	ng	96
92) Benzo(a,h,i)perylene	17.37	276	373894	10.950	ng	99

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

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