

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020819\
 Data File : BF112473.D
 Acq On : 8 Feb 2019 21:25
 Operator : JU/SJ
 Sample : K1385-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 98-SB-01-2-4

Manual Integrations
 APPROVED

Sohil
 2/11/2019 10:00:47 AM

Quant Time: Feb 09 05:58:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	153168	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	579525	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	267889	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	443830	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	291822	20.00	ng	0.00
87) Perylene-d12	15.47	264	299580	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	902427	125.14	ng	0.00
7) Phenol-d6	6.49	99	1141888	113.96	ng	0.00
23) Nitrobenzene-d5	7.40	82	740383	73.61	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	247229	79.29	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	1403366	74.09	ng	-0.02
79) Terphenyl-d14	12.94	244	1060509	68.45	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	23371	2.126	ng	90
50) Dimethylphthalate	9.58	163	111121	5.339	ng	99
71) Phenanthrene	11.38	178	98387	4.264	ng	97
75) Fluoranthene	12.57	202	201364	8.136	ng	96
78) Pyrene	12.80	202	176053	8.714	ng	95
81) Benzo(a)anthracene	13.99	228	91865	5.355	ng	99
83) Chrysene	14.03	228	81540	4.979	ng	97
86) Indeno(1,2,3-cd)pyrene	16.96	276	44488	3.429	ng	95
88) Benzo(b)fluoranthene	15.04	252	119637m	6.678	ng	
89) Benzo(k)fluoranthene	15.07	252	44714m	2.606	ng	
90) Benzo(a)pyrene	15.42	252	82178	5.098	ng	96
92) Benzo(a,h,i)perylene	17.41	276	43770	3.571	ng	98

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

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