

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113482.D
 Acq On : 1 Apr 2019 22:35
 Operator : JU/SJ
 Sample : K2182-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S8A(2-10)COMP

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:49:44 PM

Quant Time: Apr 02 03:44:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	126110	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	480448	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	224345	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	413016	20.00	ng	0.00
76) Chrysene-d12	13.85	240	369400	20.00	ng	0.00
87) Perylene-d12	15.26	264	251795	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	747614	96.93	ng	0.01
7) Phenol-d6	6.36	99	994166	101.89	ng	0.00
23) Nitrobenzene-d5	7.27	82	580481	71.71	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	265076	95.65	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1145704	78.75	ng	0.00
79) Terphenyl-d14	12.80	244	1160205	69.23	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.46	163	101251	6.090	ng	Qvalue 100
71) Phenanthrene	11.24	178	168044	8.193	ng	98
75) Fluoranthene	12.43	202	215466	9.293	ng	99
78) Pyrene	12.66	202	286916	11.208	ng	98
81) Benzo(a)anthracene	13.84	228	111223	5.260	ng	98
83) Chrysene	13.87	228	131546	6.125	ng	99
88) Benzo(b)fluoranthene	14.86	252	103721m	6.631	ng	
89) Benzo(k)fluoranthene	14.88	252	40345m	2.937	ng	
90) Benzo(a)pyrene	15.20	252	49215	3.549	ng	# 96
92) Benzo(g,h,i)perylene	17.04	276	31187	2.580	ng	# 91

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

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