

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
 Data File : BF113962.D
 Acq On : 24 Apr 2019 15:36
 Operator : JU/SJ
 Sample : K2542-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-8

Manual Integrations
 APPROVED

Sohil
 4/25/2019 5:27:02 PM

Quant Time: Apr 25 05:55:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	130003	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	493650	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	249862	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	460597	20.00	ng	0.00
76) Chrysene-d12	14.06	240	322594	20.00	ng	-0.01
87) Perylene-d12	15.56	264	413777	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	711125	93.56	ng	0.00
7) Phenol-d6	6.52	99	913883	95.83	ng	0.00
23) Nitrobenzene-d5	7.45	82	555714	69.51	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	255126	83.82	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	967044	60.53	ng	0.00
79) Terphenyl-d14	13.00	244	946775	60.17	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	188799	10.448	ng	98
71) Phenanthrene	11.43	178	61029	2.637	ng	98
75) Fluoranthene	12.62	202	137699	5.452	ng	97
78) Pyrene	12.85	202	111184	4.928	ng	97
81) Benzo(a)anthracene	14.04	228	52366	2.755	ng	96
83) Chrysene	14.08	228	59142	3.195	ng	98
86) Indeno(1,2,3-cd)pyrene	17.04	276	51666	2.947	ng	96
88) Benzo(b)fluoranthene	15.13	252	94216m	3.599	ng	
90) Benzo(a)pyrene	15.50	252	62279	2.750	ng	99
92) Benzo(g,h,i)perylene	17.49	276	48959	2.282	ng	99

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

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