

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041919\
 Data File : BF113857.D
 Acq On : 19 Apr 2019 14:41
 Operator : JU/SJ
 Sample : K2193-07 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-05-041719-A

Manual Integrations
 APPROVED

Sohil
 4/22/2019 7:26:35 PM

Quant Time: Apr 20 00:06:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	171612	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	560578	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	253556	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	520242	20.00	ng	0.00
76) Chrysene-d12	14.05	240	551758	20.00	ng	0.00
87) Perylene-d12	15.53	264	473477	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	480684	46.57	ng	0.00
7) Phenol-d6	6.52	99	588181	45.70	ng	0.00
23) Nitrobenzene-d5	7.45	82	331650	36.03	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	145090	51.79	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	543466	32.69	ng	0.00
79) Terphenyl-d14	13.00	244	750658	28.53	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	54332	3.062	ng	98
71) Phenanthrene	11.44	178	223956	8.452	ng	98
72) Anthracene	11.49	178	55239	2.069	ng	98
75) Fluoranthene	12.63	202	271687	9.675	ng	95
78) Pyrene	12.85	202	239562	6.161	ng	99
81) Benzo(a)anthracene	14.04	228	135109	4.189	ng	95
83) Chrysene	14.07	228	113917	3.558	ng	95
88) Benzo(b)fluoranthene	15.09	252	117272m	3.973	ng	
90) Benzo(a)pyrene	15.46	252	78897	3.023	ng	# 89

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

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