

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040419\
 Data File : BF113609.D
 Acq On : 5 Apr 2019 3:45
 Operator : JU/SJ
 Sample : K2240-08
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-(0-6)

Manual Integrations
 APPROVED

Sohil
 4/6/2019 12:39:22 AM

Quant Time: Apr 05 07:34:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	119196	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	412009	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	171547	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	325321	20.00	ng	0.00
76) Chrysene-d12	13.83	240	204249	20.00	ng	-0.02
87) Perylene-d12	15.23	264	167027	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	674538	96.50	ng	0.00
7) Phenol-d6	6.34	99	857002	99.48	ng	0.00
23) Nitrobenzene-d5	7.25	82	407754	57.97	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	189044	91.74	ng	-0.01
45) 2-Fluorobiphenyl	9.04	172	809285	75.95	ng	-0.01
79) Terphenyl-d14	12.77	244	717281	71.50	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.43	163	111756	8.838	ng	99
75) Fluoranthene	12.40	202	54737	3.160	ng	97
78) Pyrene	12.63	202	48122	3.096	ng	98
81) Benzo(a)anthracene	13.82	228	24489	2.078	ng	93
83) Chrysene	13.85	228	25906m	2.169	ng	
88) Benzo(b)fluoranthene	14.84	252	32842m	3.212	ng	
90) Benzo(a)pyrene	15.18	252	20363	2.241	ng	# 71

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

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