

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032719\
 Data File : BF113339.D
 Acq On : 27 Mar 2019 19:10
 Operator : JU/SJ
 Sample : K2097-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-22

Manual Integrations
 APPROVED

Sohil
 3/28/2019 8:41:16 AM

Quant Time: Mar 28 04:12:33 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	6.70	152	168896	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	619244	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	284841	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	506659	20.00	ng	0.00
76) Chrysene-d12	13.83	240	386590	20.00	ng	-0.01
87) Perylene-d12	15.24	264	428385	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	906872	87.80	ng	0.00
7) Phenol-d6	6.35	99	1197522	91.64	ng	0.00
23) Nitrobenzene-d5	7.26	82	776461	74.42	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	273796	77.82	ng	-0.01
45) 2-Fluorobiphenyl	9.06	172	1284293	66.92	ng	0.00
79) Terphenyl-d14	12.79	244	986003	56.22	ng	-0.01
Target Compounds						
10) Phenol	6.36	94	107855	7.230	ng	Qvalue # 64
50) Dimethylphthalate	9.45	163	120917	5.728	ng	99
75) Fluoranthene	12.42	202	183017	6.434	ng	99
78) Pyrene	12.64	202	170029	6.347	ng	98
81) Benzo(a)anthracene	13.82	228	76828	3.472	ng	95
83) Chrysene	13.86	228	175664	7.815	ng	98
86) Indeno(1,2,3-cd)pyrene	16.58	276	86459	4.482	ng	97
88) Benzo(b)fluoranthene	14.84	252	284397m	10.687	ng	
89) Benzo(k)fluoranthene	14.86	252	86178m	3.688	ng	
90) Benzo(a)pyrene	15.17	252	82481	3.496	ng	# 97
92) Benzo(g,h,i)perylene	16.99	276	78792	3.831	ng	99

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

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