

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032919\
 Data File : BF113414.D
 Acq On : 29 Mar 2019 15:32
 Operator : JU/SJ
 Sample : K2050-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-3

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:30 PM

Quant Time: Mar 30 01:30:16 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	158561	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	595868	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	291346	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	601643	20.00	ng	0.00
76) Chrysene-d12	13.84	240	493742	20.00	ng	0.00
87) Perylene-d12	15.24	264	497378	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	1086949	112.09	ng	0.00
7) Phenol-d6	6.35	99	1444214	117.73	ng	0.00
23) Nitrobenzene-d5	7.27	82	878853	87.54	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	387603	107.70	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1306921	66.46	ng	0.00
79) Terphenyl-d14	12.79	244	1274264	56.89	ng	0.00
Target Compounds						
10) Phenol	6.36	94	94177	6.725	ng	Qvalue # 69
32) Benzoic acid	7.77	122	111319	17.361	ng	99
50) Dimethylphthalate	9.45	163	166226	7.699	ng	99
71) Phenanthrene	11.23	178	73390	2.456	ng	96
75) Fluoranthene	12.42	202	488996	14.478	ng	99
78) Pyrene	12.64	202	451290	13.190	ng	99
81) Benzo(a)anthracene	13.83	228	191797	6.786	ng	98
83) Chrysene	13.86	228	305808	10.653	ng	98
86) Indeno(1,2,3-cd)pyrene	16.59	276	184926	7.506	ng	96
88) Benzo(b)fluoranthene	14.85	252	521298m	16.872	ng	
89) Benzo(k)fluoranthene	14.87	252	167072m	6.157	ng	
90) Benzo(a)pyrene	15.19	252	203670	7.436	ng	100
91) Dibenzo(a,h)anthracene	16.60	278	55480	2.342	ng	# 81
92) Benzo(g,h,i)perylene	17.00	276	169171	7.084	ng	97

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

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