

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022719\
 Data File : BF112740.D
 Acq On : 28 Feb 2019 1:13
 Operator : JU/SJ
 Sample : K1595-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-CBF-B11

Manual Integrations
 APPROVED

Sohil
 2/28/2019 12:47:28 PM

Quant Time: Feb 28 04:15:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 27 16:08:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	209590	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	763858	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	322313	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	628790	20.00	ng	0.00
76) Chrysene-d12	13.88	240	503398	20.00	ng	0.00
87) Perylene-d12	15.29	264	388979	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1197078	88.85	ng	0.00
7) Phenol-d6	6.37	99	1533499	90.60	ng	0.00
23) Nitrobenzene-d5	7.30	82	882628	69.14	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	342331	100.05	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1280310	65.08	ng	0.00
79) Terphenyl-d14	12.83	244	1284750	49.47	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.39	94	64708	3.276	ng	88
50) Dimethylphthalate	9.49	163	116325	4.895	ng	99
52) Acenaphthene	9.80	154	45316	2.224	ng	97
55) Dibenzofuran	9.97	168	62703	2.117	ng	97
58) Fluorene	10.32	166	64370	3.062	ng	98
71) Phenanthrene	11.27	178	1608797	51.424	ng	99
72) Anthracene	11.32	178	422096	12.889	ng	98
73) Carbazole	11.47	167	223811	7.006	ng	98
75) Fluoranthene	12.46	202	2129668	63.538	ng	97
78) Pyrene	12.69	202	1963243	46.262	ng	99
81) Benzo(a)anthracene	13.87	228	969740	27.795	ng	98
83) Chrysene	13.90	228	921898	26.976	ng	97
84) Bis(2-ethylhexyl)phthalate	13.87	149	55723	2.536	ng	97
86) Indeno(1,2,3-cd)pyrene	16.64	276	319825m	10.977	ng	
88) Benzo(b)fluoranthene	14.89	252	896772m	35.642	ng	
89) Benzo(k)fluoranthene	14.91	252	267821m	11.752	ng	
90) Benzo(a)pyrene	15.23	252	550387	24.731	ng	99
91) Dibenzo(a,h)anthracene	16.65	278	86935m	4.578	ng	
92) Benzo(a,h,i)perylene	17.04	276	300093	15.737	ng	98

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

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