

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
 Data File : BF114758.D
 Acq On : 1 Jun 2019 15:45
 Operator : HP/JU
 Sample : K3105-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 APPLE-TRACK

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:51:24 AM

Quant Time: Jun 03 03:54:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	392255	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1465105	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	700122	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1254520	20.00	ng	0.00
76) Chrysene-d12	13.82	240	750307	20.00	ng	0.00
87) Perylene-d12	15.21	264	1002827	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	2049163	91.72	ng	0.01
7) Phenol-d6	6.33	99	2501478	92.91	ng	0.00
23) Nitrobenzene-d5	7.25	82	1513843	64.46	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	597082	89.50	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	2193786	52.80	ng	0.00
79) Terphenyl-d14	12.77	244	1734924	43.47	ng	0.00

Target Compounds

						Qvalue
50) Dimethylphthalate	9.44	163	466643	9.565	ng	99
52) Acenaphthene	9.75	154	91232	2.259	ng	95
71) Phenanthrene	11.22	178	3003557	46.934	ng	98
72) Anthracene	11.27	178	808608	12.484	ng	97
75) Fluoranthene	12.41	202	4228819	63.817	ng	96
78) Pyrene	12.63	202	4241411	66.825	ng	96
81) Benzo(a)anthracene	13.81	228	2707432	46.956	ng	97
83) Chrysene	13.84	228	2374379	43.102	ng	98
86) Indeno(1,2,3-cd)pyrene	16.54	276	1427785	25.878	ng	95
88) Benzo(b)fluoranthene	14.83	252	3011274m	47.675	ng	
89) Benzo(k)fluoranthene	14.84	252	1013302m	18.458	ng	
90) Benzo(a)pyrene	15.16	252	2375472	42.811	ng	96
91) Dibenzo(a,h)anthracene	16.54	278	398106	8.119	ng	99
92) Benzo(g,h,i)perylene	16.93	276	1245703	26.177	ng	99

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

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