

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053119\
 Data File : BF114731.D
 Acq On : 1 Jun 2019 2:20
 Operator : HP/JU
 Sample : K2966-06RX
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FESB-18(2.5-3.0)RX

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:49:50 AM

Quant Time: Jun 01 03:44:43 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	6.70	152	452568	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1641914	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	747631	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1238974	20.00	ng	0.00
76) Chrysene-d12	13.82	240	792420	20.00	ng	0.00
87) Perylene-d12	15.21	264	760967	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	2532122	98.23	ng	0.02
7) Phenol-d6	6.34	99	3090300	99.49	ng	0.01
23) Nitrobenzene-d5	7.26	82	1957042	74.36	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	688929	96.71	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	2928424	70.93	ng	0.00
79) Terphenyl-d14	12.78	244	2269150	53.84	ng	0.00

Target Compounds

						Qvalue
10) Phenol	6.35	94	119808	3.193	ng	93
49) Acenaphthylene	9.58	152	217962	3.099	ng	96
50) Dimethylphthalate	9.44	163	516452	9.913	ng	99
52) Acenaphthene	9.75	154	112523	2.609	ng	98
71) Phenanthrene	11.22	178	2018630	31.939	ng	99
72) Anthracene	11.27	178	471115	7.365	ng	96
73) Carbazole	11.42	167	206067	3.665	ng	97
75) Fluoranthene	12.40	202	3111104	47.538	ng	99
78) Pyrene	12.63	202	2978564	44.435	ng	100
81) Benzo(a)anthracene	13.81	228	1628097	26.736	ng	97
83) Chrysene	13.84	228	1526313	26.234	ng	97
86) Indeno(1,2,3-cd)pyrene	16.53	276	714129	12.255	ng	94
88) Benzo(b)fluoranthene	14.82	252	1942910m	40.538	ng	
89) Benzo(k)fluoranthene	14.84	252	467005m	11.210	ng	
90) Benzo(a)pyrene	15.15	252	1241166	29.478	ng	97
91) Dibenzo(a,h)anthracene	16.54	278	196966	5.294	ng	92
92) Benzo(g,h,i)perylene	16.93	276	650935	18.026	ng	99

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

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