

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060519\
 Data File : BG041079.D
 Acq On : 5 Jun 2019 19:31
 Operator : HP/JU
 Sample : K3076-01RX
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FESB-07(5-5.5)RX

Manual Integrations
 APPROVED

mohammad
 6/6/2019 3:42:36 PM

Quant Time: Jun 06 08:49:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 18:39:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	2119	20.00	ng	-0.01
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.95
39) Acenaphthene-d10	14.73	164	59	20.00	ng	-0.02
64) Phenanthrene-d10	17.48	188	365	20.00	ng	-0.02
76) Chrysene-d12	21.76	240	3657	20.00	ng	-0.03
87) Perylene-d12	25.10	264	10190	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	327347	2785.63	ng	0.00
7) Phenol-d6	7.25	99	469084	2584.69	ng	-0.01
23) Nitrobenzene-d5	9.29	82	320126	0.00	ng	-0.02
42) 2,4,6-Tribromophenol	16.23	330	224707	256545.27	ng	-0.02
45) 2-Fluorobiphenyl	13.37	172	619689	151257.34	ng	-0.02
79) Terphenyl-d14	20.08	244	903013	5240.65	ng	-0.02
Target Compounds						
10) Phenol	7.29	94	4287	22.031	ng	Qvalue 87
50) Dimethylphthalate	14.18	163	116177	23261.594	ng	96
71) Phenanthrene	17.52	178	29558	1554.679	ng	98
72) Anthracene	17.62	178	5904	314.859	ng	99
75) Fluoranthene	19.53	202	87799	3738.777	ng	98
77) Benzo(a)pyrene	20.08	184	14597	167.322	ng	# 93
78) Pyrene	19.89	202	82074	345.241	ng	96
81) Benzo(a)anthracene	21.74	228	47320	194.386	ng	97
83) Chrysene	21.81	228	50121	215.153	ng	92
84) Bis(2-ethylhexyl)phthalate	21.61	149	12843	98.615	ng	# 94
85) Di-n-octyl phthalate	22.85	149	12200	56.248	ng	# 72
86) Indeno(1,2,3-cd)pyrene	28.91	276	65980	231.213	ng	# 59
88) Benzo(b)fluoranthene	24.02	252	89676m	145.093	ng	
89) Benzo(k)fluoranthene	24.09	252	28638	46.824	ng	98
90) Benzo(a)pyrene	24.93	252	58853	99.806	ng	93
91) Dibenzo(a,h)anthracene	28.95	278	25372m	43.061	ng	
92) Benzo(a,h,i)perylene	30.11	276	59272	103.545	ng	# 95

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

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