

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052219\
 Data File : BF114510.D
 Acq On : 23 May 2019 3:25
 Operator : JU/SJ
 Sample : K2642-05 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 Client Sampled :
 OK-01-052119

Manual Integrations
 APPROVED

mohammad
 5/24/2019 8:49:35 AM

Quant Time: May 23 05:22:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 04:47:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	202417	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	817530	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	368609	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	654969	20.00	ng	0.00
76) Chrysene-d12	14.00	240	464046	20.00	ng	0.00
87) Perylene-d12	15.47	264	405908	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	449758	35.76	ng	-0.01
7) Phenol-d6	6.47	99	587655	39.50	ng	-0.01
23) Nitrobenzene-d5	7.39	82	331869	22.78	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	115932	30.42	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	606171	24.88	ng	0.00
79) Terphenyl-d14	12.93	244	521995	23.19	ng	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.13	128	220438	5.772	ng	96
37) 2-Methylnaphthalene	8.83	142	116849	4.742	ng	99
38) 1-Methylnaphthalene	8.92	142	103615	4.466	ng	100
50) Dimethylphthalate	9.58	163	74304	2.746	ng	99
52) Acenaphthene	9.90	154	236950	10.594	ng	99
55) Dibenzofuran	10.07	168	207073	6.314	ng	95
58) Fluorene	10.42	166	329470	13.005	ng	98
71) Phenanthrene	11.38	178	1826989	57.335	ng	100
72) Anthracene	11.43	178	531654	16.611	ng	98
73) Carbazole	11.59	167	162439	5.322	ng	98
75) Fluoranthene	12.57	202	2170365	63.249	ng	98
78) Pyrene	12.80	202	2013647	53.941	ng	99
81) Benzo(a)anthracene	13.99	228	910499	30.336	ng	99
83) Chrysene	14.03	228	902414m	30.671	ng	
86) Indeno(1,2,3-cd)pyrene	16.95	276	287475	11.055	ng	96
88) Benzo(b)fluoranthene	15.04	252	904086m	34.766	ng	
89) Benzo(k)fluoranthene	15.07	252	258786m	10.833	ng	
90) Benzo(a)pyrene	15.41	252	612784	26.775	ng	99
91) Dibenzo(a,h)anthracene	16.95	278	72125	3.793	ng	# 85
92) Benzo(a,h,i)perylene	17.40	276	291753	15.308	ng	97

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

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