

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051319\
 Data File : BF114247.D
 Acq On : 14 May 2019 00:06
 Operator : JU/SJ
 Sample : K2642-03DL 20X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-050919DL

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:03:34 PM

Quant Time: May 14 08:14:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	242276	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	979749	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	459640	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	814962	20.00	ng	0.01
76) Chrysene-d12	14.02	240	580762	20.00	ng	0.00
87) Perylene-d12	15.51	264	571552	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	54241	3.60	ng	0.01
7) Phenol-d6	6.50	99	77474	4.35	ng	0.01
23) Nitrobenzene-d5	7.42	82	41342	2.37	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	15540	3.27	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	90536	2.98	ng	0.00
79) Terphenyl-d14	12.96	244	77590	2.75	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.16	128	445614	9.737	ng	98
37) 2-Methylnaphthalene	8.85	142	185607	6.286	ng	95
38) 1-Methylnaphthalene	8.95	142	145025	5.215	ng	97
52) Acenaphthene	9.92	154	279212	10.011	ng	99
55) Dibenzofuran	10.09	168	268742	6.571	ng	96
58) Fluorene	10.44	166	392679	12.430	ng	98
71) Phenanthrene	11.40	178	1948545	49.145	ng	99
72) Anthracene	11.45	178	538946	13.533	ng	98
73) Carbazole	11.61	167	242506	6.386	ng	96
75) Fluoranthene	12.60	202	1797853	42.107	ng	99
78) Pyrene	12.83	202	1614235	34.552	ng	99
81) Benzo(a)anthracene	14.01	228	720037	19.169	ng	99
83) Chrysene	14.05	228	662823	18.000	ng	98
86) Indeno(1,2,3-cd)pyrene	16.99	276	241446	7.419	ng	96
88) Benzo(b)fluoranthene	15.07	252	712601m	19.461	ng	
89) Benzo(k)fluoranthene	15.10	252	257231m	7.647	ng	
90) Benzo(a)pyrene	15.44	252	519711	16.127	ng	99
91) Dibenzo(a,h)anthracene	16.99	278	60085	2.244	ng	# 81
92) Benzo(a,h,i)perylene	17.44	276	236068	8.797	ng	99

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

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