

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071219\
 Data File : BF115528.D
 Acq On : 12 Jul 2019 16:09
 Operator : HP/JU
 Sample : K3726-01DL 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PX-1_070919DL

Manual Integrations
 APPROVED

mohammad
 7/15/2019 2:57:23 PM

Quant Time: Jul 12 16:48:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	156660	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	621624	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	305164	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	548177	20.00	ng	0.00
76) Chrysene-d12	14.04	240	304162	20.00	ng	0.00
87) Perylene-d12	15.52	264	369288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	187847	19.61	ng	-0.01
7) Phenol-d6	6.50	99	270260	22.24	ng	-0.02
23) Nitrobenzene-d5	7.43	82	158234	14.80	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	63075	17.71	ng	-0.01
45) 2-Fluorobiphenyl	9.24	172	344701	19.77	ng	-0.01
79) Terphenyl-d14	12.99	244	296482	17.99	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.19	128	67044	2.227	ng	97
37) 2-Methylnaphthalene	8.87	142	72532	3.635	ng	97
38) 1-Methylnaphthalene	8.97	142	91590	4.832	ng	99
49) Acenaphthylene	9.78	152	167584	5.693	ng	97
50) Dimethylphthalate	9.63	163	76154	3.317	ng	97
52) Acenaphthene	9.95	154	91680	4.824	ng	98
55) Dibenzofuran	10.12	168	125003	4.632	ng	# 96
58) Fluorene	10.47	166	303226	15.717	ng	99
71) Phenanthrene	11.43	178	1603115	57.052	ng	98
72) Anthracene	11.48	178	441966	15.220	ng	99
73) Carbazole	11.63	167	92447	3.630	ng	98
75) Fluoranthene	12.62	202	1185708	39.932	ng	99
78) Pvrene	12.85	202	1382255	53.639	ng	98
81) Benzo(a)anthracene	14.03	228	510376	25.470	ng	96
83) Chrysene	14.07	228	525697m	26.134	ng	
86) Indeno(1,2,3-cd)pyrene	16.99	276	350954	20.536	ng	96
88) Benzo(b)fluoranthene	15.09	252	659874m	27.750	ng	
89) Benzo(k)fluoranthene	15.11	252	200277m	9.447	ng	
90) Benzo(a)pvrene	15.46	252	536944	26.272	ng	98
91) Dibenzo(a,h)anthracene	17.00	278	102987	5.740	ng	93
92) Benzo(g,h,i)perylene	17.44	276	324808	18.151	ng	100

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

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