

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042919\
 Data File : BF114028.D
 Acq On : 29 Apr 2019 17:44
 Operator : JU/SJ
 Sample : K2564-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-2_042419

Manual Integrations
 APPROVED

Sohil
 4/30/2019 6:42:22 PM

Quant Time: Apr 30 06:23:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	127917	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	494298	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	252721	20.00	ng	-0.01
64) Phenanthrene-d10	11.41	188	463006	20.00	ng	0.00
76) Chrysene-d12	14.04	240	322562	20.00	ng	-0.02
87) Perylene-d12	15.53	264	393719	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	591358	79.07	ng	0.00
7) Phenol-d6	6.52	99	772955	82.38	ng	0.00
23) Nitrobenzene-d5	7.45	82	466281	58.24	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	246533	80.08	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	942061	58.30	ng	-0.01
79) Terphenyl-d14	12.99	244	896036	56.95	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	119310	6.528	ng	98
71) Phenanthrene	11.43	178	225860	9.707	ng	98
72) Anthracene	11.48	178	64370	2.739	ng	97
75) Fluoranthene	12.62	202	342357	13.486	ng	95
78) Pyrene	12.84	202	302295	13.401	ng	97
81) Benzo(a)anthracene	14.03	228	181350	9.542	ng	99
83) Chrysene	14.07	228	151843	8.204	ng	96
86) Indeno(1,2,3-cd)pyrene	17.01	276	97767	5.576	ng	93
88) Benzo(b)fluoranthene	15.10	252	205972m	8.269	ng	
89) Benzo(k)fluoranthene	15.12	252	69079m	3.224	ng	
90) Benzo(a)pyrene	15.47	252	157981	7.332	ng	99
92) Benzo(g,h,i)perylene	17.45	276	89129	4.367	ng	96

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

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