

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041918\
 Data File : BF104636.D
 Acq On : 19 Apr 2018 18:58
 Operator : JU/SJ
 Sample : J2387-07RX
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20180244-AMEND-COMP-ARX

Manual Integrations
 APPROVED

Sohil
 4/20/2018 3:05:23 PM

Quant Time: Apr 20 01:26:52 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 19 14:44:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	160374	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	651477	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	283500	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	427771	20.00	ng	0.00
75) Chrysene-d12	13.99	240	262955	20.00	ng	0.00
86) Perylene-d12	15.46	264	256614	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	116699	11.13	ng	0.00
7) Phenol-d6	6.45	99	885848	68.74	ng	0.00
23) Nitrobenzene-d5	7.37	82	1000518	100.28	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	13427	4.57	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1333472	74.31	ng	0.00
78) Terphenyl-d14	12.93	244	963688	83.55	ng	0.00
Target Compounds						
10) Phenol	6.45	94	36968	2.704	ng	Qvalue # 77
31) Naphthalene	8.11	128	76429	2.356	ng	100
37) 2-Methylnaphthalene	8.80	142	46961	2.238	ng	98
48) Acenaphthylene	9.70	152	484814	16.426	ng	99
49) Dimethylphthalate	9.56	163	134179	6.002	ng	99
51) Acenaphthene	9.87	154	80940	4.495	ng	97
57) Fluorene	10.39	166	102071	5.588	ng	# 96
70) Phenanthrene	11.36	178	686685	28.825	ng	100
71) Anthracene	11.41	178	527464	21.782	ng	100
74) Fluoranthene	12.56	202	984944	39.572	ng	98
77) Pyrene	12.79	202	1380532	70.829	ng	99
80) Benzo(a)anthracene	13.98	228	910179m	57.939	ng	
82) Chrysene	14.02	228	892812	55.331	ng	98
83) Bis(2-ethylhexyl)phthalate	13.96	149	110894	9.389	ng	98
85) Indeno(1,2,3-cd)pyrene	16.91	276	330343	24.100	ng	94
87) Benzo(b)fluoranthene	15.03	252	777833m	47.993	ng	
88) Benzo(k)fluoranthene	15.06	252	304942m	19.929	ng	
89) Benzo(a)pyrene	15.40	252	860052	59.308	ng	97
90) Dibenzo(a,h)anthracene	16.92	278	113365m	9.518	ng	
91) Benzo(a,h,i)perylene	17.35	276	336881	29.195	ng	97

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

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