

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
 Data File : BF086466.D
 Acq On : 28 Apr 2016 11:23
 Operator : UM/SJ
 Sample : H2702-02RE
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-COMP-B1ARE

Manual Integrations
 APPROVED

sohil
 4/28/2016 2:14:12 PM

Quant Time: Apr 28 12:40:43 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	241148	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	884765	20.00	ng	-0.01
38) Acenaphthene-d10	9.85	164	342941	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	564830	20.00	ng	0.00
75) Chrysene-d12	13.96	240	589133	20.00	ng	0.00
86) Perylene-d12	15.37	264	411831	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	811801	58.07	ng	0.01
7) Phenol-d6	6.44	99	1100406	56.39	ng	0.00
23) Nitrobenzene-d5	7.37	82	616520	37.56	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	162239	44.86	ng	-0.01
44) 2-Fluorobiphenyl	9.17	172	929644	40.90	ng	0.00
78) Terphenyl-d14	12.91	244	753319	28.17	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.56	163	156280	6.05	ng	100
70) Phenanthrene	11.35	178	154635	5.22	ng	99
74) Fluoranthene	12.53	202	329398	11.01	ng	95
77) Pyrene	12.76	202	337301	7.95	ng	100
80) Benzo(a)anthracene	13.95	228	174280	5.55	ng	97
82) Chrysene	13.99	228	176784m	5.18	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	57639	2.28	ng	93
87) Benzo(b)fluoranthene	14.97	252	156806m	6.47	ng	
88) Benzo(k)fluoranthene	14.98	252	74202m	2.94	ng	
89) Benzo(a)pyrene	15.31	252	118428	5.18	ng	97
91) Benzo(g,h,i)perylene	17.13	276	59119	3.15	ng	97

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

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