

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041717\
 Data File : BF094466.D
 Acq On : 17 Apr 2017 22:59
 Operator : SJ/MA
 Sample : I2708-07 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Z9-BASE

Manual Integrations
 APPROVED

mohammad
 4/18/2017 4:55:49 PM

Quant Time: Apr 18 03:56:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.76	152	145400	20.00	ng	0.00
21) Naphthalene-d8	7.26	136	605342	20.00	ng	0.00
38) Acenaphthene-d10	9.39	164	281114	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	485129	20.00	ng	0.00
75) Chrysene-d12	14.46	240	368351	20.00	ng	-0.01
86) Perylene-d12	16.08	264	252544	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.30	112	277416	31.65	ng	0.00
7) Phenol-d6	5.37	99	352592	34.13	ng	-0.02
23) Nitrobenzene-d5	6.40	82	211890	21.61	ng	-0.01
41) 2,4,6-Tribromophenol	10.36	330	65953	29.99	ng	-0.01
44) 2-Fluorobiphenyl	8.58	172	378964	19.83	ng	0.00
78) Terphenyl-d14	13.19	244	282976	20.53	ng	-0.01
Target Compounds						
20) 3+4-Methylphenols	6.26	107	26444	2.48	ng	Qvalue # 58
49) Dimethylphthalate	9.08	163	134877	6.44	ng	99
70) Phenanthrene	11.24	178	60644	2.22	ng	95
74) Fluoranthene	12.70	202	105814	3.74	ng	99
77) Pyrene	12.97	202	102064	3.97	ng	98
82) Chrysene	14.49	228	41893	2.14	ng	95
87) Benzo(b)fluoranthene	15.68	252	40221m	2.60	ng	
89) Benzo(a)pyrene	16.02	252	28949	2.01	ng	97

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

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