

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032718\
 Data File : BF103987.D
 Acq On : 27 Mar 2018 22:58
 Operator : JU/SJ
 Sample : J2071-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7-DB

Manual Integrations
 APPROVED

Sohil
 3/28/2018 4:54:39 PM

Quant Time: Mar 28 06:43:28 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	147625	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	587196	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	252072	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	390757	20.00	ng	0.00
75) Chrysene-d12	14.16	240	267075	20.00	ng	-0.01
86) Perylene-d12	15.70	264	231486	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	778846	80.25	ng	0.02
7) Phenol-d6	6.60	99	975454	82.15	ng	0.00
23) Nitrobenzene-d5	7.53	82	591748	70.28	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	179103	82.64	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	913889	57.83	ng	0.00
78) Terphenyl-d14	13.09	244	626628	54.46	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.72	163	145404	7.561	ng	99
70) Phenanthrene	11.53	178	236447	11.062	ng	98
71) Anthracene	11.58	178	53502	2.464	ng	98
74) Fluoranthene	12.73	202	588837	26.739	ng	97
77) Pyrene	12.96	202	532471	27.148	ng	99
80) Benzo(a)anthracene	14.14	228	331929	19.634	ng	99
82) Chrysene	14.19	228	348972	21.654	ng	98
85) Indeno(1,2,3-cd)pyrene	17.28	276	152386	10.828	ng	# 91
87) Benzo(b)fluoranthene	15.24	252	460416m	31.482	ng	
88) Benzo(k)fluoranthene	15.27	252	110432m	7.855	ng	
89) Benzo(a)pyrene	15.63	252	252276	19.019	ng	97
90) Dibenzo(a,h)anthracene	17.29	278	43088m	3.751	ng	
91) Benzo(a,h,i)perylene	17.77	276	140393	12.564	ng	95

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

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