

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080718\
 Data File : BM016241.D
 Acq On : 08 Aug 2018 05:56
 Operator : SJ/JU
 Sample : J4313-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C-5

Manual Integrations
 APPROVED

Sohil
 8/8/2018 5:35:15 PM

Quant Time: Aug 08 07:22:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 07 16:15:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	47044	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	201148	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	112701	20.00	ng	-0.01
63) Phenanthrene-d10	17.50	188	258157	20.00	ng	0.00
75) Chrysene-d12	21.64	240	269332	20.00	ng	0.00
86) Perylene-d12	23.97	264	237032	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	234278	82.72	ng	0.00
7) Phenol-d6	7.30	99	308430	83.39	ng	0.00
23) Nitrobenzene-d5	9.32	82	231023	53.42	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	104036	72.78	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	466471	50.36	ng	0.00
78) Terphenyl-d14	20.09	244	832561	64.71	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.22	163	42707	4.330	ng	Qvalue # 99
70) Phenanthrene	17.55	178	67988	4.704	ng	97
74) Fluoranthene	19.54	202	104718	6.495	ng	99
77) Pyrene	19.90	202	95065	5.890	ng	96
80) Benzo(a)anthracene	21.62	228	59113	3.564	ng	98
82) Chrysene	21.67	228	57024m	3.584	ng	
87) Benzo(b)fluoranthene	23.27	252	60769	4.354	ng	# 97
89) Benzo(a)pyrene	23.87	252	42356	3.157	ng	96
91) Benzo(g,h,i)perylene	27.09	276	27661	2.104	ng	94

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

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