

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042619\
 Data File : BG040628.D
 Acq On : 26 Apr 2019 14:19
 Operator : JU/SJ
 Sample : K2582-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSSI-0419-S-DH-6-T

Quant Time: Apr 29 06:49:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 05:35:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	52720	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	225961	20.00	ng	0.00
39) Acenaphthene-d10	14.79	164	167335	20.00	ng	0.00
64) Phenanthrene-d10	17.53	188	413559	20.00	ng	0.00
76) Chrysene-d12	21.82	240	405314	20.00	ng	0.00
87) Perylene-d12	25.19	264	434732	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	302662	101.20	ng	0.00
7) Phenol-d6	7.29	99	484188	105.15	ng	0.00
23) Nitrobenzene-d5	9.33	82	335377	68.58	ng	0.00
42) 2,4,6-Tribromophenol	16.27	330	234012	101.74	ng	0.00
45) 2-Fluorobiphenyl	13.41	172	677275	58.45	ng	0.00
79) Terphenyl-d14	20.13	244	1149310	62.82	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.23	163	76591	5.471	ng	99
74) Di-n-butylphthalate	18.47	149	73556	2.987	ng	100
75) Fluoranthene	19.57	202	73614	2.652	ng	97
78) Pyrene	19.93	202	72139	2.840	ng	98
88) Benzo(b)fluoranthene	24.12	252	69471	2.615	ng	99

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

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