

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061719\
 Data File : BG041276.D
 Acq On : 17 Jun 2019 17:09
 Operator : HP/JU
 Sample : K3156-03RE
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HD-01-061219

Quant Time: Jun 17 17:46:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 14:58:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	72061	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	278859	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	183090	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	427640	20.00	ng	0.00
76) Chrysene-d12	21.74	240	444304	20.00	ng	0.00
87) Perylene-d12	25.05	264	496016	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	314898	80.29	ng	0.00
7) Phenol-d6	7.24	99	462640	81.38	ng	0.00
23) Nitrobenzene-d5	9.26	82	336413	50.74	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	205237	73.02	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	682751	50.68	ng	0.00
79) Terphenyl-d14	20.06	244	1023569	45.66	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.16	163	134340	8.246	ng	99
75) Fluoranthene	19.51	202	112559	3.715	ng	97
78) Pyrene	19.87	202	98758	3.296	ng	99
81) Benzo(a)anthracene	21.71	228	63195	2.101	ng	95
83) Chrysene	21.78	228	71370	2.467	ng	98
88) Benzo(b)fluoranthene	23.98	252	101688	3.308	ng	# 92
90) Benzo(a)pyrene	24.89	252	69267	2.386	ng	# 95

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

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