

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041019\
 Data File : BN005127.D
 Acq On : 10 Apr 2019 14:43
 Operator : JU/SJ
 Sample : K2343-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TP-7-S

Quant Time: Apr 11 08:14:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 27 05:27:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	8469	20.00	ng	-0.02
21) Naphthalene-d8	10.43	136	34723	20.00	ng	-0.03
39) Acenaphthene-d10	14.30	164	19725	20.00	ng	-0.03
64) Phenanthrene-d10	17.05	188	42278	20.00	ng	-0.02
76) Chrysene-d12	21.25	240	38299	20.00	ng	-0.02
87) Perylene-d12	23.48	264	37140	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.25	112	57171	116.71	ng	-0.02
7) Phenol-d6	6.83	99	74803	115.13	ng	-0.02
23) Nitrobenzene-d5	8.82	82	41225	71.82	ng	-0.03
42) 2,4,6-Tribromophenol	15.80	330	36196	113.78	ng	-0.02
45) 2-Fluorobiphenyl	12.92	172	116007	77.93	ng	-0.02
79) Terphenyl-d14	19.69	244	164123	82.50	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
50) Dimethylphthalate	13.76	163	9201	5.678	ng	# 95
71) Phenanthrene	17.09	178	9681	4.075	ng	98
75) Fluoranthene	19.12	202	13220	4.878	ng	98
78) Pyrene	19.48	202	11348	4.437	ng	100
81) Benzo(a)anthracene	21.23	228	6301	2.641	ng	99
83) Chrysene	21.29	228	5221	2.291	ng	97
88) Benzo(b)fluoranthene	22.81	252	7065	3.090	ng	97
90) Benzo(a)pyrene	23.38	252	5100	2.438	ng	97

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

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