

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP080620\
 Data File : BP002926.D
 Acq On : 06 Aug 2020 18:40
 Operator : CG/JU
 Sample : L3548-07RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CRT-004RE

Quant Time: Aug 06 19:21:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 15:30:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	452404	20.00	ng	0.00
21) Naphthalene-d8	10.49	136	1834011	20.00	ng	0.00
39) Acenaphthene-d10	14.35	164	1201463	20.00	ng	-0.01
64) Phenanthrene-d10	17.10	188	2731669	20.00	ng	-0.01
76) Chrysene-d12	21.19	240	3161832	20.00	ng	-0.01
86) Perylene-d12	23.45	264	2846653	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	2963591	100.26	ng	0.00
7) Phenol-d6	6.89	99	3861951	94.10	ng	0.00
23) Nitrobenzene-d5	8.85	82	2259823	66.74	ng	-0.01
42) 2,4,6-Tribromophenol	15.84	330	1886661	126.34	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	5953358	67.70	ng	-0.01
79) Terphenyl-d14	19.67	244	9336654	59.12	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	13.80	163	632347	6.729	ng	100
71) Phenanthrene	17.14	178	403285	2.518	ng	98
75) Fluoranthene	19.12	202	1120014	6.069	ng	98
78) Pyrene	19.46	202	1047263	4.738	ng	97
80) Butylbenzylphthalate	20.36	149	8833	3.150	ng	# 94
81) Benzo(a)anthracene	21.17	228	570624	2.650	ng	100
83) Chrysene	21.22	228	651234	3.206	ng	98
88) Benzo(b)fluoranthene	22.77	252	872492	4.343	ng	99
90) Benzo(a)pyrene	23.35	252	486225	2.679	ng	100

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

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