

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082819\
 Data File : BG042552.D
 Acq On : 29 Aug 2019 3:57
 Operator : HP/JU
 Sample : K4543-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-3-S

Quant Time: Aug 29 08:03:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	33586	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	126224	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	93054	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	241092	20.00	ng	0.00
76) Chrysene-d12	21.68	240	297885	20.00	ng	-0.01
87) Perylene-d12	24.92	264	356734	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	152235	91.80	ng	0.00
7) Phenol-d6	7.16	99	199355	89.00	ng	-0.01
23) Nitrobenzene-d5	9.16	82	111831	61.22	ng	-0.01
42) 2,4,6-Tribromophenol	16.15	330	135445	102.41	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	351183	63.83	ng	0.00
79) Terphenyl-d14	20.02	244	781595	56.72	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.19	94	12794	5.617	ng	94
50) Dimethylphthalate	14.11	163	44225	6.587	ng	99
71) Phenanthrene	17.45	178	29720	2.482	ng	95
75) Fluoranthene	19.46	202	86561	5.923	ng	95
78) Pyrene	19.83	202	91686	5.020	ng	96
81) Benzo(a)anthracene	21.67	228	57653	3.182	ng	96
83) Chrysene	21.73	228	58582	3.352	ng	97
88) Benzo(b)fluoranthene	23.89	252	69013	3.519	ng	100
90) Benzo(a)pyrene	24.77	252	45377	2.438	ng	99

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

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