

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090519\
 Data File : BG042745.D
 Acq On : 6 Sep 2019 8:08
 Operator : HP/JU
 Sample : K4692-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-24-COMP

Quant Time: Sep 06 08:53:45 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	7.99	152	38631	20.00	ng	-0.01
21) Naphthalene-d8	10.82	136	136044	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	89856	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	195843	20.00	ng	-0.01
76) Chrysene-d12	21.68	240	228000	20.00	ng	-0.01
87) Perylene-d12	24.93	264	273235	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.56	112	256048	134.24	ng	0.00
7) Phenol-d6	7.17	99	323899	125.71	ng	0.00
23) Nitrobenzene-d5	9.17	82	185605	94.28	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	181364	142.01	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	510631	96.11	ng	0.00
79) Terphenyl-d14	20.02	244	857288	81.29	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
50) Dimethylphthalate	14.10	163	63033	9.722	ng	98
71) Phenanthrene	17.45	178	45773	4.705	ng	98
75) Fluoranthene	19.46	202	84565	7.123	ng	93
78) Pyrene	19.83	202	88438	6.327	ng	99
81) Benzo(a)anthracene	21.67	228	54457	3.926	ng	94
83) Chrysene	21.73	228	50498	3.775	ng	98
88) Benzo(b)fluoranthene	23.90	252	68467	4.558	ng	97
90) Benzo(a)pyrene	24.77	252	52704	3.697	ng	99
92) Benzo(g,h,i)perylene	29.79	276	33437	2.316	ng	97

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

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