

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090119\
 Data File : BM022476.D
 Acq On : 01 Sep 2019 14:30
 Operator : HP/JU
 Sample : K4619-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP1-COMP

Quant Time: Sep 02 04:07:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 02 03:54:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	11758	20.00	ng	0.00
21) Naphthalene-d8	10.28	136	44065	20.00	ng	0.00
39) Acenaphthene-d10	14.17	164	29271	20.00	ng	0.00
64) Phenanthrene-d10	16.94	188	75963	20.00	ng	0.00
76) Chrysene-d12	21.16	240	131379	20.00	ng	0.00
87) Perylene-d12	23.34	264	349171	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	71489	106.72	ng	0.00
7) Phenol-d6	6.71	99	91476	102.12	ng	0.00
23) Nitrobenzene-d5	8.68	82	64438	64.05	ng	0.00
42) 2,4,6-Tribromophenol	15.69	330	54880	91.62	ng	0.00
45) 2-Fluorobiphenyl	12.77	172	142948	55.32	ng	0.00
79) Terphenyl-d14	19.59	244	459561	46.16	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.65	163	13135	5.159	ng	99
60) Diethylphthalate	15.02	149	6511	2.448	ng	95
71) Phenanthrene	16.99	178	17592	3.973	ng	99
73) Carbazole	17.38	167	8379	2.327	ng	99
75) Fluoranthene	19.02	202	95325	16.166	ng	97
78) Pyrene	19.38	202	81810	9.137	ng	99
81) Benzo(a)anthracene	21.14	228	76602	8.123	ng	99
83) Chrysene	21.19	228	60563	6.748	ng	96

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

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