

Data Path : V:\HPCHEM1\BNA M\DATA\BM042016\
 Data File : BM004995.D
 Acq On : 20 Apr 2016 17:50
 Operator : UM/SJ
 Sample : H2621-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H2TR3

Quant Time: Apr 21 10:20:35 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM041416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 15 22:59:24 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	35189	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	159473	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.45	164	98612	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	222500	20.00	ng/ul	0.00
75) Chrysene-d12	21.38	240	262136	20.00	ng/ul	0.00
83) Perylene-d12	23.67	264	245934	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.31	96	1757	2.67	ng/uL	0.00
5) Phenol-d5	6.97	99	36104	14.37	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	22678	15.74	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	30089	14.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	28759	13.86	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	12183	11.53	ng/ul	0.00
22) 2-Nitrophenol-d4	9.69	143	13416	11.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	27794	12.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.74	131	32464	13.35	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	89151	13.29	ng/ul	0.00
46) Acenaphthylene-d8	14.14	160	111011	13.49	ng/ul	0.00
51) 4-Nitrophenol-d4	14.65	143	6000	5.24	ng/ul	0.00
57) Fluorene-d10	15.44	176	84448	14.57	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.56	200	7074	7.35	ng/ul	0.00
70) Anthracene-d10	17.29	188	125798	14.23	ng/ul	0.00
76) Pyrene-d10	19.59	212	149082	13.65	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.52	264	141207	14.91	ng/ul	0.00

Target Compounds						Ovalue
44) Dimethylphthalate	13.90	163	23555	3.53	ng/ul	98

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

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