

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071916\
 Data File : BF089123.D
 Acq On : 19 Jul 2016 17:40
 Operator : UM/SJ
 Sample : H4025-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HSR-WC-60-071116

Quant Time: Jul 20 03:37:40 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 19 15:28:17 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	48337	20.00	ng	0.00
21) Naphthalene-d8	7.91	136	200789	20.00	ng	0.00
38) Acenaphthene-d10	9.66	164	88022	20.00	ng	0.00
63) Phenanthrene-d10	11.13	188	139423	20.00	ng	0.00
75) Chrysene-d12	13.75	240	97880	20.00	ng	-0.01
86) Perylene-d12	15.10	264	79470	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	376800	116.83	ng	0.01
7) Phenol-d6	6.28	99	532955	127.89	ng	0.00
23) Nitrobenzene-d5	7.19	82	336615	87.85	ng	-0.01
41) 2,4,6-Tribromophenol	10.44	330	88904	108.77	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	634068	113.79	ng	0.00
78) Terphenyl-d14	12.71	244	374061	85.12	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	2.19	88	7187	4.49	ng	Qvalue # 87
3) Pyridine	2.81	79	105306	22.70	ng	94
4) n-Nitrosodimethylamine	2.74	42	9972	5.63	ng	95
31) Naphthalene	7.93	128	33143	3.35	ng	99
33) 4-Chloroaniline	7.99	127	33970	7.71	ng	96
35) Caprolactam	8.31	113	5911	6.53	ng	93
76) Benzidine	12.46	184	105590	21.34	ng	97

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

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