

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091719\
 Data File : BP000269.D
 Acq On : 17 Sep 2019 13:48
 Operator : HP/JU
 Sample : K4897-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-200052

Quant Time: Sep 18 12:40:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 16 18:59:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	295230	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1011127	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	561517	20.00	ng	0.01
64) Phenanthrene-d10	11.36	188	913480	20.00	ng	0.04
76) Chrysene-d12	14.13	240	96326	20.00	ng	0.13
87) Perylene-d12	15.63	264	235177	20.00	ng	0.15
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1786508	104.47	ng	0.01
7) Phenol-d6	6.43	99	2265839	108.09	ng	0.00
23) Nitrobenzene-d5	7.35	82	1086689	69.34	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	689523	123.80	ng	0.02
45) 2-Fluorobiphenyl	9.15	172	3030008	97.12	ng	0.00
79) Terphenyl-d14	13.02	244	800392	174.31	ng	0.08
Target Compounds						
10) Phenol	6.44	94	126061	5.293	ng	98
37) 2-Methylnaphthalene	8.79	142	325183	10.158	ng	99
38) 1-Methylnaphthalene	8.89	142	178402	5.949	ng	99
50) Dimethylphthalate	9.55	163	227393	5.944	ng	99
66) n-Nitrosodiphenylamine	10.50	169	77818	2.830	ng	98
71) Phenanthrene	11.40	178	158793	3.469	ng	# 59
86) Indeno(1,2,3-cd)pyrene	17.10	276	40797	5.602	ng	# 67

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

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