

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080621\
 Data File : BP006555.D
 Acq On : 07 Aug 2021 01:03
 Operator : CG/JU
 Sample : M3298-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1924

Quant Time: Aug 07 05:41:38 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:41:30 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.751	152	215442	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	892128	20.000	ng	#-0.01
39) Acenaphthene-d10	14.381	164	508910	20.000	ng	-0.01
64) Phenanthrene-d10	17.139	188	958523	20.000	ng	# 0.00
76) Chrysene-d12	21.233	240	933101	20.000	ng	# 0.00
86) Perylene-d12	23.557	264	997400	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	948072	67.591	ng	0.00
7) Phenol-d6	6.934	99	732764	36.776	ng	0.00
23) Nitrobenzene-d5	8.910	82	1811340	93.716	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	774860	145.687	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	3037005	89.284	ng	0.00
79) Terphenyl-d14	19.710	244	4287028	92.062	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	8.522	107	172906	10.069	ng	Qvalue 90

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

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