

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062321\
 Data File : BP006038.D
 Acq On : 23 Jun 2021 17:01
 Operator : CG/JU
 Sample : M2789-21
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1ST-FLOOR

Quant Time: Jun 24 01:16:54 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 22 17:52:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	84565	20.000	ng	0.00
21) Naphthalene-d8	10.722	136	326436	20.000	ng	0.00
39) Acenaphthene-d10	14.546	164	201305	20.000	ng	0.00
64) Phenanthrene-d10	17.316	188	354645	20.000	ng	# 0.02
76) Chrysene-d12	21.398	240	441369	20.000	ng	# 0.03
86) Perylene-d12	23.810	264	496102	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	344732	65.413	ng	0.00
7) Phenol-d6	7.093	99	424602	62.160	ng	0.00
23) Nitrobenzene-d5	9.081	82	275005	41.303	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	240896	69.733	ng	0.01
45) 2-Fluorobiphenyl	13.175	172	685987	44.748	ng	0.00
79) Terphenyl-d14	19.881	244	1052157	44.637	ng	0.04
Target Compounds						
						Qvalue
32) Benzoic acid	10.069	122	22958	11.733	ng	97
71) Phenanthrene	17.357	178	761888	35.773	ng	# 83
75) Fluoranthene	19.339	202	132374	5.118	ng	# 12
78) Pyrene	19.710	202	105429	3.422	ng	# 1
83) Chrysene	21.433	228	84430	2.736	ng	# 32
84) Bis(2-ethylhexyl)phtha...	21.292	149	597245	32.124	ng	# 88

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

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