

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120721\
 Data File : BP008262.D
 Acq On : 07 Dec 2021 21:47
 Operator : CG/JU
 Sample : M4949-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-221

Quant Time: Dec 08 01:11:11 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 03 14:39:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	58579	20.000	ng	-0.02
21) Naphthalene-d8	10.581	136	205587	20.000	ng	-0.02
39) Acenaphthene-d10	14.434	164	125893	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	271352	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	333467	20.000	ng	-0.01
86) Perylene-d12	23.698	264	403433	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	480775	132.145	ng	-0.01
7) Phenol-d6	6.969	99	596832	121.521	ng	-0.02
23) Nitrobenzene-d5	8.946	82	404509	89.472	ng	-0.02
42) 2,4,6-Tribromophenol	15.934	330	235790	146.533	ng	-0.02
45) 2-Fluorobiphenyl	13.051	172	788171	90.938	ng	-0.02
79) Terphenyl-d14	19.769	244	1266001	79.344	ng	-0.01
Target Compounds						
						Qvalue
10) Phenol	6.999	94	18433	3.652	ng	93
75) Fluoranthene	19.216	202	57523	2.980	ng	96
81) Benzo(a)anthracene	21.286	228	49167	2.225	ng	98
83) Chrysene	21.339	228	56708	2.696	ng	98
84) Bis(2-ethylhexyl)phtha...	21.216	149	56630	3.928	ng	97
88) Benzo(b)fluoranthene	22.968	252	93943	3.863	ng	99
90) Benzo(a)pyrene	23.592	252	63478	3.054	ng	# 95
92) Benzo(g,h,i)perylene	26.951	276	57252	2.328	ng	96

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

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