Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121021\
Data File : BP008357.D
Acq On : 11 Dec 2021 00:30
Operator : CG/JU
Sample : M4959-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-1S

Quant Time: Dec 11 04:02:59 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 11 03:53:01 2021
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	74114	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	226709	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	127333	20.000	ng	0.00
64) Phenanthrene-d10	17.175	188	266828	20.000	ng	0.00
76) Chrysene-d12	21.280	240	342553	20.000	ng	0.00
86) Perylene-d12	23.662	264	390191	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	540221	117.361	ng	0.00
7) Phenol-d6	6.952	99	649224	104.480	ng	0.00
23) Nitrobenzene-d5	8.922	82	421597	84.564	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	210891	129.577	ng	0.00
45) 2-Fluorobiphenyl	13.028	172	705837	80.518	ng	0.00
79) Terphenyl-d14	19.751	244	1168820	71.310	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.198	202	80541	4.244	ng	99
81) Benzo(a)anthracene	21.268	228	63323	2.789	ng	92
83) Chrysene	21.321	228	70519	3.264	ng	96
88) Benzo(b)fluoranthene	22.939	252	98683	4.195	ng	97
90) Benzo(a)pyrene	23.557	252	73559	3.659	ng	95
92) Benzo(g,h,i)perylene	26.898	276	52520	2.208	ng	99

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

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